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Publication number: **0 612 846 A1**

(12)

EUROPEAN PATENT APPLICATION

(21) Application number: **94101207.2**

(51) Int. Cl.⁵: **C12N 15/27, C07K 3/00,
C12P 21/02, C07K 13/00,
G06F 15/60**

(22) Date of filing: **27.01.94**

(83) Declaration under Rule 28(4) EPC (expert solution)

(30) Priority: **28.01.93 US 10099**

(43) Date of publication of application:
31.08.94 Bulletin 94/35

(64) Designated Contracting States:
**AT BE CH DE DK ES FR GB GR IE IT LI LU MC
NL PT SE**

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(54) **G-CSF analog compositions and methods.**

(57) Provided herein are granulocyte colony stimulating factor ("G-CSF") analogs, compositions containing such analogs, and related compositions. In another aspect, provided herein are nucleic acids encoding the present analogs or related nucleic acids, related host cells and vectors. In yet another aspect, provided herein are computer programs and apparatuses for expressing the three dimensional structure of G-CSF and analogs thereof. In another aspect, provided herein are methods for rationally designing G-CSF analogs and related compositions. In yet another aspect, provided herein are methods for treatment using the present G-CSF analogs.

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Field of the Invention

This invention relates to granulocyte colony stimulating factor ("G-CSF") analogs, compositions containing such analogs, and related compositions. In another aspect, the present invention relates to nucleic acids encoding the present analogs or related nucleic acids, related host cells and vectors. In another aspect, the invention relates to computer programs and apparatuses for expressing the three dimensional structure of G-CSF and analogs thereof. In another aspect, the invention relates to methods for rationally designing G-CSF analogs and related compositions. In yet another aspect, the present invention relates to methods for treatment using the present G-CSF analogs.

Background

Hematopoiesis is controlled by two systems: the cells within the bone marrow microenvironment and growth factors. The growth factors, also called colony stimulating factors, stimulate committed progenitor cells to proliferate and to form colonies of differentiating blood cells. One of these factors is granulocyte colony stimulating factor, herein called G-CSF, which preferentially stimulates the growth and development of neutrophils, indicating a potential use in neutropenic states. Welte et al., PNAS-USA 82: 1526-1530 (1985); Souza et al., Science 232: 61-65 (1986) and Gabrilove, J. Seminars in Hematology 26: (2) 1-14 (1989).

In humans, endogenous G-CSF is detectable in blood plasma. Jones et al., Bailliere's Clinical Hematology 2 (1): 83-111 (1989). G-CSF is produced by fibroblasts, macrophages, T cells trophoblasts, endothelial cells and epithelial cells and is the expression product of a single copy gene comprised of four exons and five introns located on chromosome seventeen. Transcription of this locus produces a mRNA species which is differentially processed, resulting in two forms of G-CSF mRNA, one version coding for a protein of 177 amino acids, the other coding for a protein of 174 amino acids, Nagata et al., EMBO J 5: 575-581 (1986), and the form comprised of 174 amino acids has been found to have the greatest specific *in vivo* biological activity. G-CSF is species cross-reactive, such that when human G-CSF is administered to another mammal such as a mouse, canine or monkey, sustained neutrophil leukocytosis is elicited. Moore et al., PNAS-USA 84: 7134-7138 (1987).

Human G-CSF can be obtained and purified from a number of sources. Natural human G-CSF (nhG-CSF) can be isolated from the supernatants of cultured human tumor cell lines. The development of recombinant DNA technology, see, for instance, U.S. Patent 4,810,643 (Souza) incorporated herein by reference, has enabled the production of commercial scale quantities of G-CSF in glycosylated form as a product of eukaryotic host cell expression, and of G-CSF in non-glycosylated form as a product of prokaryotic host cell expression.

G-CSF has been found to be useful in the treatment of indications where an increase in neutrophils will provide benefits. For example, for cancer patients, G-CSF is beneficial as a means of selectively stimulating neutrophil production to compensate for hematopoietic deficits resulting from chemotherapy or radiation therapy. Other indications include treatment of various infectious diseases and related conditions, such as sepsis, which is typically caused by a metabolite of bacteria. G-CSF is also useful alone, or in combination with other compounds, such as other cytokines, for growth or expansion of cells in culture, for example, for bone marrow transplants.

Signal transduction, the way in which G-CSF effects cellular metabolism, is not currently thoroughly understood. G-CSF binds to a cell-surface receptor which apparently initiates the changes within particular progenitor cells, leading to cell differentiation.

Various altered G-CSF's have been reported. Generally, for design of drugs, certain changes are known to have certain structural effects. For example, deleting one cysteine could result in the unfolding of a molecule which is, in its unaltered state, is normally folded via a disulfide bridge. There are other known methods for adding, deleting or substituting amino acids in order to change the function of a protein.

Recombinant human G-CSF mutants have been prepared, but the method of preparation does not include overall structure/function relationship information. For example, the mutation and biochemical modification of Cys 18 has been reported. Kuga et al., Biochem. Biophys. Res. Comm 159: 103-111 (1989); Lu et al., Arch. Biochem. Biophys. 268: 81-92 (1989).

In U.S. Patent No. 4, 810, 643, entitled, "Production of Pluripotent Granulocyte Colony-Stimulating Factor" (as cited above), polypeptide analogs and peptide fragments of G-CSF are disclosed generally. Specific G-CSF analogs disclosed include those with the cysteins at positions 17, 36, 42, 64, and 74 (of the 174 amino acid species or of those having 175 amino acids, the additional amino acid being an N-terminal methionine) substituted with another amino acid, (such as serine), and G-CSF with an alanine in the first (N-

terminal) position.

EP 0 335 423 entitled "Modified human G-CSF" reportedly discloses the modification of at least one amino group in a polypeptide having hG-CSF activity.

EP 0 272 703 entitled "Novel Polypeptide" reportedly discloses G-CSF derivatives having an amino acid substituted or deleted at or "in the neighborhood" of the N terminus.

EP 0 459 630, entitled "Polypeptides" reportedly discloses derivatives of naturally occurring G-CSF having at least one of the biological properties of naturally occurring G-CSF and a solution stability of at least 35% at 5 mg/ml in which the derivative has at least Cys¹⁷ of the native sequence replaced by a Ser¹⁷ residue and Asp²⁷ of the native sequence replaced by a Ser²⁷ residue.

EP 0 256 843 entitled "Expression of G-CSF and Muteins Thereof and Their Uses" reportedly discloses a modified DNA sequence encoding G-CSF wherein the N-terminus is modified for enhanced expression of protein in recombinant host cells, without changing the amino acid sequence of the protein.

EP 0 243 153 entitled "Human G-CSF Protein Expression" reportedly discloses G-CSF to be modified by inactivating at least one yeast KEX2 protease processing site for increased yield in recombinant production using yeast.

Shaw, U.S. Patent No. 4,904,584, entitled "Site-Specific Homogeneous Modification of Polypeptides," reportedly discloses lysine altered proteins.

WO/9012874 reportedly discloses cysteine altered variants of proteins.

Australian patent application Document No. AU-A-10948/92, entitled, "Improved Activation of Recombinant Proteins" reportedly discloses the addition of amino acids to either terminus of a G-CSF molecule for the purpose of aiding in the folding of the molecule after prokaryotic expression.

Australian patent application Document No. AU-A-76380/91, entitled, "Muteins of the Granulocyte Colony Stimulating Factor (G-CSF)" reportedly discloses muteins of the granulocyte stimulating factor G-CSF in the sequence Leu-Gly-His-Ser-Leu-Gly-Ile at position 50-56 of G-CSF with 174 amino acids, and position 53 to 59 of the G-CSF with 177 amino acids, or/and at least one of the four histidine residues at positions 43, 79, 156 and 170 of the mature G-CSF with 174 amino acids or at positions 46, 82, 159, or 173 of the mature G-CSF with 177 amino acids.

GB 2 213 821, entitled "Synthetic Human Granulocyte Colony Stimulating Factor Gene" reportedly discloses a synthetic G-CSF-encoding nucleic acid sequence incorporating restriction sites to facilitate the cassette mutagenesis of selected regions, and flanking restriction sites to facilitate the incorporation of the gene into a desired expression system.

G-CSF has reportedly been crystallized to some extent, e.g., EP 344 796, and the overall structure of G-CSF has been surmised, but only on a gross level. Bazan, Immunology Today 11: 350-354 (1990); Parry et al., J. Molecular Recognition 8: 107-110 (1988). To date, there have been no reports of the overall structure of G-CSF, and no systematic studies of the relationship of the overall structure and function of the molecule, studies which are essential to the systematic design of G-CSF analogs. Accordingly, there exists a need for a method of this systematic design of G-CSF analogs, and the resultant compositions.

Summary of the Invention

The three dimensional structure of G-CSF has now been determined to the atomic level. From this three-dimensional structure, one can now forecast with substantial certainty how changes in the composition of a G-CSF molecule may result in structural changes. These structural characteristics may be correlated with biological activity to design and produce G-CSF analogs.

Although others had speculated regarding the three dimensional structure of G-CSF, Bazan, Immunology Today 11: 350-354 (1990); Parry et al., J. Molecular Recognition 8: 107-110 (1988), these speculations were of no help to those wishing to prepare G-CSF analogs either because the surmised structure was incorrect (Parry et al., *supra*) and/or because the surmised structure provided no detail correlating the constituent moieties with structure. The present determination of the three-dimensional structure to the atomic level is by far the most complete analysis to date, and provides important information to those wishing to design and prepare G-CSF analogs. For example, from the present three dimensional structural analysis, precise areas of hydrophobicity and hydrophilicity have been determined.

Relative hydrophobicity is important because it directly relates to the stability of the molecule. Generally, biological molecules, found in aqueous environments, are externally hydrophilic and internally hydrophobic; in accordance with the second law of thermodynamics provides, this is the lowest energy state and provides for stability. Although one could have speculated that G-CSF's internal core would be hydrophobic, and the outer areas would be hydrophilic, one would have had no way of knowing specific hydrophobic or hydrophilic areas. With the presently provided knowledge of areas of hydrophobic-

ity/philicity, one may forecast with substantial certainty which changes to the G-CSF molecule will affect the overall structure of the molecule.

As a general rule, one may use knowledge of the geography of the hydrophobic and hydrophilic regions to design analogs in which the overall G-CSF structure is not changed, but change does affect biological activity ("biological activity" being used here in its broadest sense to denote function). One may correlate biological activity to structure. If the structure is not changed, and the mutation has no effect on biological activity, then the mutation has no biological function. If, however, the structure is not changed and the mutation does affect biological activity, then the residue (or atom) is essential to at least one biological function. Some of the present working examples were designed to provide no change in overall structure, yet have a change in biological function.

Based on the correlation of structure to biological activity, one aspect of the present invention relates to G-CSF analogs. These analogs are molecules which have more, fewer, different or modified amino acid residues from the G-CSF amino acid sequence. The modifications may be by addition, substitution, or deletion of one or more amino acid residues. The modification may include the addition or substitution of analogs of the amino acids themselves, such as peptidomimetics or amino acids with altered moieties such as altered side groups. The G-CSF used as a basis for comparison may be of human, animal or recombinant nucleic acid-technology origin (although the working examples disclosed herein are based on the recombinant production of the 174 amino acid species of human G-CSF, having an extra N-terminus methionyl residue). The analogs may possess functions different from natural human G-CSF molecule, or may exhibit the same functions, or varying degrees of the same functions. For example, the analogs may be designed to have a higher or lower biological activity, have a longer shelf-life or a decrease in stability, be easier to formulate, or more difficult to combine with other ingredients. The analogs may have no hematopoietic activity, and may therefore be useful as an antagonist against G-CSF effect (as, for example, in the overproduction of G-CSF). From time to time herein the present analogs are referred to as proteins or peptides for convenience, but contemplated herein are other types of molecules, such as peptidomimetics or chemically modified peptides.

In another aspect, the present invention relates to related compositions containing a G-CSF analog as an active ingredient. The term, "related composition," as used herein, is meant to denote a composition which may be obtained once the identity of the G-CSF analog is ascertained (such as a G-CSF analog labeled with a detectable label, related receptor or pharmaceutical composition). Also considered a related composition are chemically modified versions of the G-CSF analog, such as those having attached at least one polyethylene glycol molecule.

For example, one may prepare a G-CSF analog to which a detectable label is attached, such as a fluorescent, chemiluminescent or radioactive molecule.

Another example is a pharmaceutical composition which may be formulated by known techniques using known materials, see, e.g., Remington's Pharmaceutical Sciences, 18th Ed. (1990, Mack Publishing Co., Easton, Pennsylvania 18042) pages 1435-1712, which are herein incorporated by reference. Generally, the formulation will depend on a variety of factors such as administration, stability, production concerns and other factors. The G-CSF analog may be administered by injection or by pulmonary administration via inhalation. Enteric dosage forms may also be available for the present G-CSF analog compositions, and therefore oral administration may be effective. G-CSF analogs may be inserted into liposomes or other microcarriers for delivery, and may be formulated in gels or other compositions for sustained release. Although preferred compositions will vary depending on the use to which the composition will be put, generally, for G-CSF analogs having at least one of the biological activities of natural G-CSF, preferred pharmaceutical compositions are those prepared for subcutaneous injection or for pulmonary administration via inhalation, although the particular formulations for each type of administration will depend on the characteristics of the analog.

Another example of related composition is a receptor for the present analog. As used herein, the term "receptor" indicates a moiety which selectively binds to the present analog molecule. For example, antibodies, or fragments thereof, or "recombinant antibodies" (see Huse et al., Science 246:1275 (1989)) may be used as receptors. Selective binding does not mean only specific binding (although binding-specific receptors are encompassed herein), but rather that the binding is not a random event. Receptors may be on the cell surface or intra- or extra-cellular, and may act to effectuate, inhibit or localize the biological activity of the present analogs. Receptor binding may also be a triggering mechanism for a cascade of activity indirectly related to the analog itself. Also contemplated herein are nucleic acids, vectors containing such nucleic acids and host cells containing such nucleic acids which encode such receptors.

Another example of a related composition is a G-CSF analog with a chemical moiety attached. Generally, chemical modification may alter biological activity or antigenicity of a protein, or may alter other

characteristics, and these factors will be taken into account by a skilled practitioner. As noted above, one example of such chemical moiety is polyethylene glycol. Modification may include the addition of one or more hydrophilic or hydrophobic polymer molecules, fatty acid molecules, or polysaccharide molecules. Examples of chemical modifiers include polyethylene glycol, alkylpolyethylene glycols, Di-poly(amino acids),
 5 polyvinylpyrrolidone, polyvinyl alcohol, pyran copolymer, acetic acid/acylation, propionic acid, palmitic acid, stearic acid, dextran, carboxymethyl cellulose, pullulan, or agarose. See, Francis, *Focus on Growth Factors* 3: 4-10 (May 1992) (published by Mediscript, Mountview Court, Friern Barnet Lane, London N20 OLD, UK). Also, chemical modification may include an additional protein or portion thereof, use of a cytotoxic agent, or an antibody. The chemical modification may also include lecithin.

10 In another aspect, the present invention relates to nucleic acids encoding such analogs. The nucleic acids may be DNAs or RNAs or derivatives thereof, and will typically be cloned and expressed on a vector, such as a phage or plasmid containing appropriate regulatory sequences. The nucleic acids may be labeled (such as using a radioactive, chemiluminescent, or fluorescent label) for diagnostic or prognostic purposes, for example. The nucleic acid sequence may be optimized for expression, such as including codons
 15 preferred for bacterial expression. The nucleic acid and its complementary strand, and modifications thereof which do not prevent encoding of the desired analog are here contemplated.

In another aspect, the present invention relates to host cells containing the above nucleic acids encoding the present analogs. Host cells may be eukaryotic or prokaryotic, and expression systems may include extra steps relating to the attachment (or prevention) of sugar groups (glycosylation), proper folding
 20 of the molecule, the addition or deletion of leader sequences or other factors incident to recombinant expression.

In another aspect the present invention relates to antisense nucleic acids which act to prevent or modify the type or amount of expression of such nucleic acid sequences. These may be prepared by known methods.

25 In another aspect of the present invention, the nucleic acids encoding a present analog may be used for gene therapy purposes, for example, by placing a vector containing the analog-encoding sequence into a recipient so the nucleic acid itself is expressed inside the recipient who is in need of the analog composition. The vector may first be placed in a carrier, such as a cell, and then the carrier placed into the recipient. Such expression may be localized or systemic. Other carriers include non-naturally occurring
 30 carriers, such as liposomes or other microcarriers or particles, which may act to mediate gene transfer into a recipient.

The present invention also provides for computer programs for the expression (such as visual display) of the G-CSF or analog three dimensional structure, and further, a computer program which expresses the identity of each constituent of a G-CSF molecule and the precise location within the overall structure of that
 35 constituent, down to the atomic level. Set forth below is one example of such program. There are many currently available computer programs for the expression of the three dimensional structure of a molecule. Generally, these programs provide for inputting of the coordinates for the three dimensional structure of a molecule (i.e., for example, a numerical assignment for each atom of a G-CSF molecule along an x, y, and z axis), means to express (such as visually display) such coordinates, means to alter such coordinates and
 40 means to express an image of a molecule having such altered coordinates. One may program crystallographic information, i.e., the coordinates of the location of the atoms of a G-CSF molecule in three dimension space, wherein such coordinates have been obtained from crystallographic analysis of said G-CSF molecule, into such programs to generate a computer program for the expression (such as visual display) of the G-CSF three dimensional structure. Also provided, therefore, is a computer program for the expression
 45 of G-CSF analog three dimensional structure. Preferred is the computer program Insight II, version 4, available from Biosym, San Diego, California, with the coordinates as set forth in FIGURE 5 input. Preferred expression means is on a Silicon Graphics 320 VGX computer, with Crystal Eyes glasses (also available from Silicon Graphics), which allows one to view the G-CSF molecule or its analog stereoscopically. Alternatively, the present G-CSF crystallographic coordinates and diffraction data are also deposited in the
 50 Protein Data Bank, Chemistry Department, Brookhaven National Laboratory, Upton, New York 119723, USA. One may use these data in preparing a different computer program for expression of the three dimensional structure of a G-CSF molecule or analog thereof. Therefore, another aspect of the present invention is a computer program for the expression of the three dimensional structure of a G-CSF molecule. Also provided is said computer program for visual display of the three dimensional structure of a G-CSF
 55 molecule; and further, said program having means for altering such visual display. Apparatus useful for expression of such computer program, particularly for the visual display of the computer image of said three dimensional structure of a G-CSF molecule or analog thereof is also therefore here provided, as well as means for preparing said computer program and apparatus.

The computer program is useful for preparation of G-CSF analogs because one may select specific sites on the G-CSF molecule for alteration and readily ascertain the effect the alteration will have on the overall structure of the G-CSF molecule. Selection of said site for alteration will depend on the desired biological characteristic of the G-CSF analog. If one were to randomly change said G-CSF molecule (r-met-hu-G-CSF) there would be 175²⁰ possible substitutions, and even more analogs having multiple changes, additions or deletions. By viewing the three dimensional structure wherein said structure is correlated with the composition of the molecule, the selection for sites of alteration is no longer a random event, but sites for alteration may be determined rationally.

As set forth above, identity of the three dimensional structure of G-CSF, including the placement of each constituent down to the atomic level has now yielded information regarding which moieties are necessary to maintain the overall structure of the G-CSF molecule. One may therefore select whether to maintain the overall structure of the G-CSF molecule when preparing a G-CSF analog of the present invention, or whether (and how) to change the overall structure of the G-CSF molecule when preparing a G-CSF analog of the present invention. Optionally, once one has prepared such analog, one may test such analog for a desired characteristic.

One may, for example, seek to maintain the overall structure possessed by a non-altered natural or recombinant G-CSF molecule. The overall structure is presented in Figures 2, 3, and 4, and is described in more detail below. Maintenance of the overall structure may ensure receptor binding, a necessary characteristic for an analog possessing the hematopoietic capabilities of natural G-CSF (if no receptor binding, signal transduction does not result from the presence of the analog). It is contemplated that one class of G-CSF analogs will possess the three dimensional core structure of a natural or recombinant (non-altered) G-CSF molecule, yet possess different characteristics, such as an increased ability to selectively stimulate neutrophils. Another class of G-CSF analogs are those with a different overall structure which diminishes the ability of a G-CSF analog molecule to bind to a G-CSF receptor, and possesses a diminished ability to selectively stimulate neutrophils as compared to non-altered natural or recombinant G-CSF.

For example, it is now known which moieties within the internal regions of the G-CSF molecule are hydrophobic, and, correspondingly, which moieties on the external portion of the G-CSF molecule are hydrophilic. Without knowledge of the overall three dimensional structure, preferably to the atomic level as provided herein, one could not forecast which alterations within this hydrophobic internal area would result in a change in the overall structural conformation of the molecule. An overall structural change could result in a functional change, such as lack of receptor binding, for example, and therefore, diminishment of biological activity as found in non-altered G-CSF. Another class of G-CSF analogs is therefore G-CSF analogs which possess the same hydrophobicity as (non-altered) natural or recombinant G-CSF. More particularly, another class of G-CSF analogs possesses the same hydrophobic moieties within the four helical bundle of its internal core as those hydrophobic moieties possessed by (non-altered) natural or recombinant G-CSF yet have a composition different from said non-altered natural or recombinant G-CSF.

Another example relates to external loops which are structures which connect the internal core (helices) of the G-CSF molecule. From the three dimensional structure -- including information regarding the spatial location of the amino acid residues -- one may forecast that certain changes in certain loops will not result in overall conformational changes. Therefore, another class of G-CSF analogs provided herein is that having an altered external loop but possessing the same overall structure as (non-altered) natural or recombinant G-CSF. More particularly, another class of G-CSF analogs provided herein are those having an altered external loop, said loop being selected from the loop present between helices A and B; between helices B and C; between helices C and D; between helices D and A, as those loops and helices are identified herein. More particularly, said loops, preferably the AB loop and/or the CD loop are altered to increase the half life of the molecule by stabilizing said loops. Such stabilization may be by connecting all or a portion of said loop(s) to a portion of an alpha helical bundle found in the core of a G-CSF (or analog) molecule. Such connection may be via beta sheet, salt bridge, disulfide bonds, hydrophobic interaction or other connecting means available to those skilled in the art, wherein such connecting means serves to stabilize said external loop or loops. For example, one may stabilize the AB or CD loops by connecting the AB loop to one of the helices within the internal region of the molecule.

The N-terminus also may be altered without change in the overall structure of a G-CSF molecule, because the N-terminus does not effect structural stability of the internal helices, and, although the external loops are preferred for modification, the same general statements apply to the N-terminus.

Additionally, such external loops may be the site(s) for chemical modification because in (non-altered) natural or recombinant G-CSF such loops are relatively flexible and tend not to interfere with receptor binding. Thus, there would be additional room for a chemical moiety to be directly attached (or indirectly

attached via another chemical moiety which serves as a chemical connecting means). The chemical moiety may be selected from a variety of moieties available for modification of one or more function of a G-CSF molecule. For example, an external loop may provide sites for the addition of one or more polymer which serves to increase serum half-life, such as a polyethylene glycol molecule. Such polyethylene glycol molecule(s) may be added wherein said loop is altered to include additional lysines which have reactive side groups to which polyethylene glycol moieties are capable of attaching. Other classes of chemical moieties may also be attached to one or more external loops, including but not limited to other biologically active molecules, such as receptors, other therapeutic proteins (such as other hematopoietic factors which would engender a hybrid molecule), or cytotoxic agents (such as diphtheria toxin). This list is of course not complete; one skilled in the art possessed of the desired chemical moiety will have the means to effect attachment of said desired moiety to the desired external loop. Therefore, another class of the present G-CSF analogs includes those with at least one alteration in an external loop wherein said alteration provides for the addition of a chemical moiety such as at least one polyethylene glycol molecule.

Deletions, such as deletions of sites recognized by proteins for degradation of the molecule, may also be effectual in the external loops. This provides alternative means for increasing half-life of a molecule otherwise having the G-CSF receptor binding and signal transduction capabilities (i.e., the ability to selectively stimulate the maturation of neutrophils). Therefore, another class of the present G-CSF analogs includes those with at least one alteration in an external loop wherein said alteration decreases the turnover of said analog by proteases. Preferred loops for such alterations are the AB loop and the CD loop. One may prepare an abbreviated G-CSF molecule by deleting a portion of the amino acid residues found in the external loops (identified in more detail below), said abbreviated G-CSF molecule may have additional advantages in preparation or in biological function.

Another example relates to the relative charges between amino acid residues which are in proximity to each other. As noted above, the G-CSF molecule contains a relatively tightly packed four helical bundle. Some of the faces on the helices face other helices. At the point (such as a residue) where a helix faces another helix, the two amino acid moieties which face each other may have the same charge, and thus tend to repel each other, which lends instability to the overall molecule. This may be eliminated by changing the charge (to an opposite charge or a neutral charge) of one or both of the amino acid moieties so that there is no repelling. Therefore, another class of G-CSF analogs includes those G-CSF analogs having been altered to modify instability due to surface interactions, such as electron charge location.

In another aspect, the present invention relates to methods for designing G-CSF analogs and related compositions and the products of those methods. The end products of the methods may be the G-CSF analogs as defined above or related compositions. For instance, the examples disclosed herein demonstrate (a) the effects of changes in the constituents (i.e., chemical moieties) of the G-CSF molecule on the G-CSF structure and (b) the effects of changes in structure on biological function. Essentially, therefore, another aspect of the present invention is a method for preparing a G-CSF analog comprising the steps of:

- (a) viewing information conveying the three dimensional structure of a G-CSF molecule wherein the chemical moieties, such as each amino acid residue or each atom of each amino acid residue, of the G-CSF molecule are correlated with said structure;
- (b) selecting from said information a site on a G-CSF molecule for alteration;
- (c) preparing a G-CSF analog molecule having such alteration; and
- (d) optionally, testing such G-CSF analog molecule for a desired characteristic.

One may use the here provided computer programs for a computer-based method for preparing a G-CSF analog. Another aspect of the present invention is therefore a computer based method for preparing a G-CSF analog comprising the steps of:

- (a) providing computer expression of the three dimensional structure of a G-CSF molecule wherein the chemical moieties, such as each amino acid residue or each atom of each amino acid residue, of the G-CSF molecule are correlated with said structure;
- (b) selecting from said computer expression a site on a G-CSF molecule for alteration;
- (c) preparing a G-CSF molecule having such alteration; and
- (d) optionally, testing such G-CSF molecule for a desired characteristic.

More specifically, the present invention provides a method for preparing a G-CSF analog comprising the steps of:

- (a) viewing the three dimensional structure of a G-CSF molecule via a computer, said computer programmed (i) to express the coordinates of a G-CSF molecule in three dimensional space, and (ii) to allow for entry of information for alteration of said G-CSF expression and viewing thereof;
- (b) selecting a site on said visual image of said G-CSF molecule for alteration;
- (c) entering information for said alteration on said computer;

- (d) viewing a three dimensional structure of said altered G-CSF molecule via said computer;
- (e) optionally repeating steps (a)-(e);
- (f) preparing a G-CSF analog with said alteration; and
- (g) optionally testing said G-CSF analog for a desired characteristic.

5 In another aspect, the present invention relates to methods of using the present G-CSF analogs and related compositions and methods for the treatment or protection of mammals, either alone or in combination with other hematopoietic factors or drugs in the treatment of hematopoietic disorders. It is contemplated that one aspect of designing G-CSF analogs will be the goal of enhancing or modifying the characteristics non-modified G-CSF is known to have.

10 For example, the present analogs may possess enhanced or modified activities, so, where G-CSF is useful in the treatment of (for example) neutropenia, the present compositions and methods may also be of such use.

Another example is the modification of G-CSF for the purpose of interacting more effectively when used in combination with other factors particularly in the treatment of hematopoietic disorders. One example of such combination use is to use an early-acting hematopoietic factor (i.e., a factor which acts earlier in the hematopoiesis cascade on relatively undifferentiated cells) and either simultaneously or in seriatim use of a later-acting hematopoietic factor, such as G-CSF or analog thereof (as G-CSF acts on the CFU-GM lineage in the selective stimulation of neutrophils). The present methods and compositions may be useful in therapy involving such combinations or "cocktails" of hematopoietic factors.

20 The present compositions and methods may also be useful in the treatment of leukopenia, myelogenous leukemia, severe chronic neutropenia, aplastic anemia, glycogen storage disease, mucositis, and other bone marrow failure states. The present compositions and methods may also be useful in the treatment of hematopoietic deficits arising from chemotherapy or from radiation therapy. The success of bone marrow transplantation, or the use of peripheral blood progenitor cells for transplantation, for example, may be enhanced by application of the present compositions (proteins or nucleic acids for gene therapy) and methods. The present compositions and methods may also be useful in the treatment of infectious diseases, such in the context of wound healing, burn treatment, bacteremia, septicemia, fungal infections, endocarditis, osteomyelitis, infection related to abdominal trauma, infections not responding to antibiotics, pneumonia and the treatment of bacterial inflammation may also benefit from the application of the present compositions and methods. In addition, the present compositions and methods may be useful in the treatment of leukemia based upon a reported ability to differentiate leukemic cells. Welte et al., PNAS-USA 82: 1526-1530 (1985). Other applications include the treatment of individuals with tumors, using the present compositions and methods, optionally in the presence of receptors (such as antibodies) which bind to the tumor cells. For review articles on therapeutic applications, see Lieshke and Burgess, N.Engl.J.Med. 327: 28-34 and 99-106 (1992) both of which are herein incorporated by reference.

35 The present compositions and methods may also be useful to act as intermediaries in the production of other moieties; for example, G-CSF has been reported to influence the production of other hematopoietic factors and this function (if ascertained) may be enhanced or modified via the present compositions and/or methods.

40 The compositions related to the present G-CSF analogs, such as receptors, may be useful to act as an antagonist which prevents the activity of G-CSF or an analog. One may obtain a composition with some or all of the activity of non-altered G-CSF or a G-CSF analog, and add one or more chemical moieties to alter one or more properties of such G-CSF or analog. With knowledge of the three dimensional conformation, one may forecast the best geographic location for such chemical modification to achieve the desired effect.

45 General objectives in chemical modification may include improved half-life (such as reduced renal, immunological or cellular clearance), altered bioactivity (such as altered enzymatic properties, dissociated bioactivities or activity in organic solvents), reduced toxicity (such as concealing toxic epitopes, compartmentalization, and selective biodistribution), altered immunoreactivity (reduced immunogenicity, reduced antigenicity or adjuvant action), or altered physical properties (such as increased solubility, improved thermal stability, improved mechanical stability, or conformational stabilization). See Francis, *Focus on Growth Factors* 3: 4-10 (May 1992) (published by Mediscript, Mountview Court, Friern Barnet Lane, London N20 0LD, UK).

50 The examples below are illustrative of the present invention and are not intended as a limitation. It is understood that variations and modifications will occur to those skilled in the art, and it is intended that the appended claims cover all such equivalent variations which come within the scope of the invention as claimed.

Detailed Description of the Drawings

FIGURE 1 is an illustration of the amino acid sequence of the 174 amino acid species of G-CSF with an additional N-terminal methionine (Seq. ID No.: 1) (Seq. ID No.: 2).

FIGURE 2 is a topology diagram of the crystalline structure of G-CSF, as well as hGH, pGH, GM-CSF, INF-B, IL-2, and IL-4. These illustrations are based on inspection of cited references. The length of secondary structural elements are drawn in proportion to the number of residues. A, B, C, and D helices are labeled according to the scheme used herein for G-CSF. For INF- β , the original labeling of helices is indicated in parentheses.

FIGURE 3 is a "ribbon diagram" of the three dimensional structure of G-CSF. Helix A is amino acid residues 11-39 (numbered according to Figure 1, above), helix B is amino acid residues 72-91, helix C is amino acid residues 100-123, and helix D is amino acid residues 143-173. The relatively short 3^{10} helix is at amino acid residues 45-48, and the alpha helix is at amino acid residues 48-53. Residues 93-95 form almost one turn of a left handed helix.

FIGURE 4 is a "barrel diagram" of the three dimensional structure of G-CSF. Shown in various shades of gray are the overall cylinders and their orientations for the three dimensional structure of G-CSF. The numbers indicate amino acid residue position according to FIGURE 1 above.

FIGURE 5 is a list of the coordinates used to generate a computer-aided visual image of the three-dimensional structure of G-CSF. The coordinates are set forth below. The columns correspond to separate

field:

- (i) Field 1 (from the left hand side) is the atom,
- (ii) Field 2 is the assigned atom number,
- (iii) Field 3 is the atom name (according to the periodic table standard nomenclature, with CB being carbon atom Beta, CG is Carbon atom Gamma, etc.);
- (iv) Field 4 is the residue type (according to three letter nomenclature for amino acids as found in, e.g., Stryer, Biochemistry, 3d Ed., W.H. Freeman and Company, N.Y. 1988, inside back cover);
- (v) Fields 5-7 are the x-axis, y-axis and z-axis positions of the atom;
- (vi) Field 8 (often a "1.00") designates occupancy at that position;
- (vii) Field 9 designates the B-factor;
- (viii) Field 10 designates the molecule designation. Three molecules (designated a, b, and c) of G-CSF crystallized together as a unit. The designation a, b, or c indicates which coordinates are from which molecule. The number after the letter (1, 2, or 3) indicates the assigned amino acid residue position, with molecule A having assigned positions 10-175, molecule B having assigned positions 210-375, and molecule C having assigned positions 410-575. These positions were so designated so that there would be no overlap among the three molecules which crystallized together. (The "W" designation indicates water).

FIGURE 6 is a schematic representation of the strategy involved in refining the crystallization matrix for parameters involved in crystallization. The crystallization matrix corresponds to the final concentration of the components (salts, buffers and precipitants) of the crystallization solutions in the wells of a 24 well tissue culture plate. These concentrations are produced by pipetting the appropriate volume of stock solutions into the wells of the microtiter plate. To design the matrix, the crystallographer decides on an upper and lower concentration of the component. These upper and lower concentrations can be pipetted along either the rows (e.g., A1-A6, B1-B6, C1-C6 or D1-D6) or along the entire tray (A1-D6). The former method is useful for checking reproducibility of crystal growth of a single component along a limited number of wells, whereas the later method is more useful in initial screening. The results of several stages of refinement of the crystallization matrix are illustrated by a representation of three plates. The increase in shading in the wells indicates a positive crystallization result which, in the final stages, would be X-ray quality crystals but in the initial stages could be oil droplets, granular precipitates or small crystals approximately less than 0.05 mm in size. Part A represents an initial screen of one parameter in which the range of concentration between the first well (A1) and last well (D6) is large and the concentration increase between wells is calculated as $(\text{concentration A1}) - (\text{concentration D6}) / 23$. Part B represents that in later stages of the crystallization matrix refinement of the concentration spread between A1 and D6 would be reduced which would result in more crystals formed per plate. Part C indicates a final stage of matrix refinement in which quality crystals are found in most wells of the plate.

Detailed Description of the Invention

The present invention grows out of the discovery of the three dimensional structure of G-CSF. This three dimensional structure has been expressed via computer program for stereoscopic viewing. By viewing this stereoscopically, structure-function relationships identified and G-CSF analogs have been designed and made.

The Overall Three Dimensional Structure of G-CSF

The G-CSF used to ascertain the structure was a non-glycosylated 174 amino acid species having an extra N-terminal methionine residue incident to bacterial expression. The DNA and amino acid sequence of this G-CSF are illustrated in FIGURE 1.

Overall, the three dimensional structure of G-CSF is predominantly helical, with 103 of the 175 residues forming a 4-alpha-helical bundle. The only other secondary structure is found in the loop between the first two long helices where a 4 residue 3^{10} helix is immediately followed by a 6 residue alpha helix. As shown in FIGURE 2, the overall structure has been compared with the structure reported for other proteins: growth hormone (Abdel-Meguid et al., PNAS-USA 84: 6434 (1987) and Vos et al., Science 255: 305-312 (1992)), granulocyte macrophage colony stimulating factor (Diederichs et al., Science 254: 1779-1782 (1991)), interferon- β (Senda et al., EMBO J. 11: 3193-3201 (1992)), interleukin-2 (McKay Science 257: 1673-1677 (1992)) and interleukin-4 (Powers et al., Science 256: 1673-1677 (1992), and Smith et al., J. Mol. Biol. 224: 899-904 (1992)). Structural similarity among these growth factors occurs despite the absence of similarity in their amino acid sequences.

Presently, the structural information was correlation of G-CSF biochemistry, and this can be summarized as follows (with sequence position 1 being at the N-terminus):

Sequence Position	Description of Structure	Analysis
1-10	Extended chain	Deletion causes no loss of biological activity
Cys 18	Partially buried	Reactive with DTNB and Thimersosol but not with iodo-acetate
34	Alternative splice site	Insertion reduces biological activity
20-47 (inclusive)	Helix A, first disulfide and portion of AB helix	Predicted receptor binding region based on neutralizing antibody data
20, 23, 24	Helix A	Single alanine mutation of residue(s) reduces biological activity. Predicted receptor binding (Site B).
165-175 (inclusive)	Carboxy terminus	Deletion reduces biological activity

This biochemical information, having been gleaned from antibody binding studies, see Layton et al., Biochemistry 266: 23815-23823 (1991), was superimposed on the three-dimensional structure in order to design G-CSF analogs. The design, preparation, and testing of these G-CSF analogs is described in Example 1 below.

EXAMPLE 1

This Example describes the preparation of crystalline G-CSF, the visualization of the three dimensional structure of recombinant human G-CSF via computer-generated image, the preparation of analogs, using site-directed mutagenesis or nucleic acid amplification methods, the biological assays and HPLC analysis used to analyze the G-CSF analogs, and the resulting determination of overall structure/function relationships. All cited publications are herein incorporated by reference.

A. Use of Automated Crystallization

The need for a three-dimensional structure of recombinant human granulocyte colony stimulating factor (r-hu-G-CSF), and the availability of large quantities of the purified protein, led to methods of crystal growth by incomplete factorial sampling and seeding. Starting with the implementation of incomplete factorial

crystallization described by Jancarik and Kim, *J. Appl. Crystallogr.* **24**: 409 (1991) solution conditions that yielded oil droplets and birefringence aggregates were ascertained. Also, software and hardware of an automated pipetting system were modified to produce some 400 different crystallization conditions per day. Weber, *J. Appl. Crystallogr.* **20**: 366-373 (1987). This procedure led to a crystallization solution which produced r-hu-G-CSF crystals.

The size, reproducibility and quality of the crystals was improved by a seeding method in which the number of "nucleation initiating units" was estimated by serial dilution of a seeding solution. These methods yielded reproducible growth of 2.0 mm r-hu-G-CSF crystals. The space group of these crystals is P2₁2₁2₁ with cell dimensions of a = 90 Å, b = 110 Å and c = 49 Å, and they diffract to a resolution of 2.0 Å.

1. Overall Methodology

To search for the crystallizing conditions of a new protein, Carter and Carter, *J. Biol. Chem.* **254**: 122219-12223 (1979) proposed the incomplete factorial method. They suggested that a sampling of a large number of randomly selected, but generally probable, crystallizing conditions may lead to a successful combination of reagents that produce protein crystallization. This idea was implemented by Jancarik and Kim, *J. Appl. Crystallogr.* **24**: 409(1991), who described 32 solutions for the initial crystallization trials which cover a range of pH, salts and precipitants. Here we describe an extension of their implementation to an expanded set of 70 solutions. To minimize the human effort and error of solution preparation, the method has been programmed for an automatic pipetting machine.

Following Weber's method of successive automated grid searching (SAGS), *J. Cryst. Growth* **90**: 318-324(1988), the robotic system was used to generate a series of solutions which continually refined the crystallization conditions of temperature, pH, salts and precipitant. Once a solution that could reproducibly grow crystals was determined, a seeding technique which greatly improved the quality of the crystals was developed. When these methods were combined, hundreds of diffraction quality crystals (crystals diffracting to at least about 2.5 Angstroms, preferably having at least portions diffracting to below 2 Angstroms, and more preferably, approximately 1 Angstrom) were produced in a few days.

Generally, the method for crystallization, which may be used with any protein one desires to crystallize, comprises the steps of:

- (a) combining aqueous aliquots of the desired protein with either (i) aliquots of a salt solution, each aliquot having a different concentration of salt; or (ii) aliquots of a precipitant solution, each aliquot having a different concentration of precipitant, optionally wherein each combined aliquot is combined in the presence of a range of pH;
- (b) observing said combined aliquots for precrystalline formations, and selecting said salt or precipitant combination and said pH which is efficacious in producing precrystalline forms, or, if no precrystalline forms are so produced, increasing the protein starting concentration of said aqueous aliquots of protein;
- (c) after said salt or said precipitant concentration is selected, repeating step (a) with said previously unselected solution in the presence of said selected concentration; and
- (d) repeating step (b) and step (a) until a crystal of desired quality is obtained.

The above method may optionally be automated, which provides vast savings in time and labor. Preferred protein starting concentrations are between 10mg/ml and 20mg/ml, however this starting concentration will vary with the protein (the G-CSF below was analyzed using 33mg/ml). A preferred range of salt solution to begin analysis with is (NaCl) of 0-2.5M. A preferred precipitant is polyethylene glycol 8000, however, other precipitants include organic solvents (such as ethanol), polyethylene glycol molecules having a molecular weight in the range of 500-20,000, and other precipitants known to those skilled in the art. The preferred pH range is pH 4.5, 5.0, 5.5, 6.0, 6.5, 7.0, 7.5, 8.0, 8.5, and 9.0. Precrystallization forms include oils, birefringement precipitants, small crystals (< approximately 0.05 mm), medium crystals (approximately 0.5 to .5 mm) and large crystals (> approximately 0.5 mm). The preferred time for waiting to see a crystalline structure is 48 hours, although weekly observation is also preferred, and generally, after about one month, a different protein concentration is utilized (generally the protein concentration is increased). Automation is preferred, using the Accuflex system as modified. The preferred automation parameters are described below.

Generally, protein with a concentration between 10 mg/ml and 20 mg/ml was combined with a range of NaCl solutions from 0-2.5 M, and each such combination was performed (separately) in the presence of the above range of concentrations. Once a precrystallization structure is observed, that salt concentration and pH range are optimized in a separate experiment, until the desired crystal quality is achieved. Next, the precipitant concentration, in the presence of varying levels of pH is also optimized. When both are optimized, the optimal conditions are performed at once to achieve the desired result (this is diagrammed in

FIGURE 6).

a. Implementation of an automated pipetting system

5 Drops and reservoir solutions were prepared by an Accuflex pipetting system (ICN Pharmaceuticals, Costa Mesa, CA) which is controlled by a personal computer that sends ASCII codes through a standard serial interface. The pipetter samples six different solutions by means of a rotating valve and pipettes these solutions onto a plate whose translation in a x-y coordinate system can be controlled. The vertical component of the system manipulates a syringe that is capable both of dispensing and retrieving liquid.

10 The software provided with the Accuflex was based on the SAGS method as proposed by Cox and Weber, J.Appl. Crystallogr. 20: 366-373 (1987). This method involves the systematic variation of two major crystallization parameters, pH and precipitant concentration, with provision to vary two others. While building on these concepts, the software used here provided greater flexibility in the design and implementation of the crystallization solutions used in the automated grid searching strategy. As a result of this flexibility the present software also created a larger number of different solutions. This is essential for
15 the implementation of the incomplete factorial method as described in that section below.

To improve the speed and design of the automated grid searching strategy, the Accuflex pipetting system required software and hardware modifications. The hardware changes allowed the use of two different micro-titer trays, one used for hanging drop and one used for sitting drop experiments, and a
20 Plexiglas tray which held 24 additional buffer, salt and precipitant solutions. These additional solutions expanded the grid of crystallizing conditions that could be surveyed.

To utilize the hardware modifications, the pipetting software was written in two subroutines; one subroutine allows the crystallographer to design a matrix of crystallization solutions based on the concentrations of their components and the second subroutine to translate these concentrations into the computer
25 code which pipettes the proper volumes of the solutions into the crystallization trays. The concentration matrices can be generated by either of two programs. The first program (MRF, available from Amgen, Inc., Thousand Oaks, CA) refers to a list of stock solution concentrations supplied by the crystallographer and calculates the required volume to be pipette to achieve the designated concentration. The second method, which is preferred, incorporates a spread sheet program (Lotus) which can be used to make more
30 sophisticated gradients of precipitants or pH. The concentration matrix created by either program is interpreted by the control program (SUX, a modification of the program found in the Accuflex pipetter originally and available from Amgen, Inc., Thousand Oaks, CA) and the wells are filled accordingly.

b. Implementation of the Incomplete Factorial Method

35 The convenience of the modified pipetting system for preparing diverse solutions improved the implementation of an expanded incomplete factorial method. The development of a new set of crystallization solutions having "random" components was generated using the program INFAC, Carter et al., J.Cryst. Growth 90: 60-73(1988) which produced a list containing 96 random combinations of one factor from three
40 variables. Combinations of calcium and phosphate which immediately precipitated were eliminated, leaving 70 distinct combinations of precipitants, salts and buffers. These combinations were prepared using the automated pipetter and incubated for 1 week. The mixtures were inspected and solutions which formed precipitants were prepared again with lower concentrations of their components. This was repeated until all wells were clear of precipitant.

c. Crystallization of r-hu-G-CSF

45 Several different crystallization strategies were used to find a solution which produced x-ray quality crystals. These strategies included the use of the incomplete factorial method, refinement of the crystallization conditions using successive automated grid searches (SAGS), implementation of a seeding technique and development of a crystal production procedure which yielded hundreds of quality crystals overnight. Unless otherwise noted the screening and production of r-hu-G-CSF crystals utilized the hanging drop vapor diffusion method. Afinsen et al., Physical principles of protein crystallization. In: Eisenberg (ed.), Advances in Protein Chemistry 41: 1-33 (1991).

55 The initial screening for crystallization conditions of r-hu-G-CSF used the Jancarik and Kim, J.Appl.Crystallogr. 24: 409(1991) incomplete factorial method which resulted in several solutions that produced "precrySTALLIZATION" results. These results included birefringent precipitants, oils and very small crystals (< .05 mm). These precrySTALLIZATIONS solutions then served as the starting points for systematic

screening.

The screening process required the development of crystallization matrices. These matrices corresponded to the concentration of the components in the crystallization solutions and were created using the IBM-PC based spread sheet Lotus™ and implemented with the modified Accuflex pipetting system.

5 The strategy in designing the matrices was to vary one crystallization condition (such as salt concentration) while holding the other conditions such as pH, and precipitant concentration constant. At the start of screening, the concentration range of the varied condition was large but the concentration was successively refined until all wells in the micro-titer tray produced the same crystallization result. These results were scored as follows: crystals, birefringent precipitate, granular precipitate, oil droplets and amorphous

10 mass. If the concentration of a crystallization parameter did not produce at least a precipitant, the concentration of that parameter was increased until a precipitant formed. After each tray was produced, it was left undisturbed for at least two days and then inspected for crystal growth. After this initial screening, the trays were then inspected on a weekly basis.

From this screening process, two independent solutions with the same pH and precipitant but differing

15 in salts (MgCl, LiSO₄) were identified which produced small (0.1 x 0.05 x 0.05 mm) crystals. Based on these results, a new series of concentration matrices were produced which varied MgCl with respect to LiSO₄ while keeping the other crystallization parameters constant. This series of experiments resulted in identification of a solution which produced diffraction quality crystals (> approximately 0.5 mm) in about three weeks. To find this crystallization growth solution (100 mM Mes pH 5.8, 380 mM MgCl₂, 220 mM

20 LiSO₄ and 8% PEG 8k) approximately 8,000 conditions had been screened which consumed about 300 mg of protein.

The size of the crystals depended on the number of crystals forming per drop. Typically 3 to 5 crystals would be formed with average size of (1.0 x 0.7 x 0.7 mm). Two morphologies which had an identical space group (P2₁2₁2₁) and unit cell dimensions a = 90.2, b = 110.2, c = 49.5 were obtained depending on whether

25 or not seeding (see below) was implemented. Without seeding, the r-hu-G-CSF crystals had one long flat surface and rounded edges.

When seeding was employed, crystals with sharp faces were observed in the drop within 4 to 6 hours (0.05 by 0.05 by 0.05 mm). Within 24 hours, crystals had grown to (0.7 by 0.7 by 0.7 mm) and continued to grow beyond 2 mm depending on the number of crystals forming in the drop.

30

d. Seeding and determination of nucleation initiation sites.

The presently provided method for seeding crystals establishes the number of nucleation initiation units in each individual well used (here, after the optimum conditions for growing crystals had been determined).

35 The method here is advantageous in that the number of "seeds" affects the quality of the crystals, and this in turn affects the degree of resolution. The present seeding here also provides advantages in that with seeding, G-CSF crystal grows in a period of about 3 days, whereas without seeding, the growth takes approximately three weeks.

In one series of production growth (see methods), showers of small but well defined crystals were

40 produced overnight (<0.01 x 0.01 x 0.01 mm). Crystallization conditions were followed as described above except that a pipette tip employed in previously had been reused. Presumably, the crystal showering effect was caused by small nucleation units which had formed in the used tip and which provided sites of nucleation for the crystals. Addition of a small amount (0.5 ul) of the drops containing the crystal showers to a new drop under standard production growth conditions resulted in a shower of crystals overnight. This

45 method was used to produce several trays of drops containing crystal showers which we termed "seed stock".

The number of nucleation initiation units (NIU) contained within the "seed stock" drops was estimated to attempt to improve the reproducibility and quality of the r-hu-GCSF crystals. To determine the number of NIU in the "seed stock", an aliquot of the drop was serially diluted along a 96 well microtiter plate. The

50 microtiter plate was prepared by adding 50 ul of a solution containing equal volumes of r-hu-G-CSF (33 mg/ml) and the crystal growth solution (described above) in each well. An aliquot (3 ul) of one of the "seed stock" drops was transferred to the first well of the microtiter plate. The solution in the well was mixed and 3 ul was then transferred to the next well along the row of the microtiter plate. Each row of the microtiter plate was similarly prepared and the tray was sealed with plastic tape. Overnight, small crystals formed in

55 the bottom of the wells of the microtiter plate and the number of crystals in the wells were correlated to the dilution of the original "seed stock". To produce large single crystals, the "seed stock" drop was appropriately diluted into fresh CGS and then an aliquot of this solution containing the NIU was transferred to a drop

Once crystallization conditions had been optimized, crystals were grown in a production method in which 3 ml each of CGS and r-hu-G-CSF (33 mg/ml) were mixed to create 5 trays (each having 24 wells). This method included the production of the refined crystallization solution in liter quantities, mixing this solution with protein and placing the protein/crystallization solution in either hanging drop or sitting drop trays. This process typically yielded 100 to 300 quality crystals (>0.5 mm) in about 5 days.

e. Experimental Methods

Materials

Crystallographic information was obtained starting with r-hu-met-G-CSF with the amino acid sequence as provided in FIGURE 1 with a specific activity of $1.0 \pm 0.6 \times 10^8$ U/mg (as measured by cell mitogenesis assay in a 10 mM acetate buffer at pH 4.0 (in Water for Injection) at a concentration of approximately 3 mg/ml solution was concentrated with an Amicon concentrator at 75 psi using a YM10 filter. The solution was typically concentrated 10 fold at 4 °C and stored for several months.

Initial Screening

Crystals suitable for X-ray analysis were obtained by vapor-diffusion equilibrium using hanging drops. For preliminary screening, 7 ul of the protein solution at 33 mg/ml (as prepared above) was mixed with an equal volume of the well solution, placed on siliconized glass plates and suspended over the well solution utilizing Linbro tissue culture plates (Flow Laboratories, McLean, Va). All of the pipetting was performed with the Accuflex pipetter, however, trays were removed from the automated pipetter after the well solutions had been created and thoroughly mixed for at least 10 minutes with a table top shaker. The Linbro trays were then returned to the pipetter which added the well and protein solutions to the siliconized cover slips. The cover slips were then inverted and sealed over 1 ml of the well solutions with silicon grease.

The components of the automated crystallization system are as follows. A PC-DOS computer system was used to design a matrix of crystallization solutions based on the concentration of their components. These matrices were produced with either MRF of the Lotus spread sheet (described above). The final product of these programs is a data file. This file contains the information required by the SUX program to pipette the appropriate volume of the stock solutions to obtain the concentrations described in the matrices. The SUX program information was passed through a serial I/O port and used to dictate to the Accuflex pipetting system the position of the valve relative to the stock solutions, the amount of solution to be retrieved, and then pipetted into the wells of the microtiter plates and the X-Y position of each well (the column/row of each well). Addition information was transmitted to the pipetter which included the Z position (height) of the syringe during filling as well as the position of a drain where the system pauses to purge the syringe between fillings of different solutions. The 24 well microtiter plate (either Linbro or Cryschem) and cover slip holder was placed on a plate which was moved in the X-Y plane. Movement of the plate allowed the pipetter to position the syringe to pipette into the wells. It also positioned the coverslips and vials and extract solutions from these sources. Prior the pipetting, the Linbro microtiter plates had a thin film of grease applied around the edges of the wells. After the crystallization solutions were prepared in the wells and before they were transferred to the cover slips, the microtiter plate was removed from the pipetting system, and solutions were allowed to mix on a table top shaker for ten minutes. After mixing, the well solution was either transferred to the cover slips (in the case of the hanging drop protocol) or transferred to the middle post in the well (in the case of the sitting drop protocol). Protein was extracted from a vial and added to the coverslip drop containing the well solution (or to the post). Plastic tape was applied to the top of the Cryschem plate to seal the wells.

Production Growth

Once conditions for crystallization had been optimized, crystal growth was performed utilizing a "production" method. The crystallization solution which contained 100 mM Mes pH 5.8, 380 mM MgCl₂, 220 mM LiSO₄, and 8% PEG 8K was made in 1 liter quantities. Utilizing an Eppendorf syringe pipetter, 1 ml aliquots of this solution were pipetted into each of the wells of the Linbro plate. A solution containing 50% of this solution and 50% G-CSF (33 mg/ml) was mixed and pipetted onto the siliconized cover slips. Typical volumes of these drops were between 50 and 100 ul and because of the large size of these drops, great care was taken in flipping the coverslips and suspending the drops over the wells.

Data Collection

The structure has been refined with X-PLOR (Bruniger, X-PLOR version 3.0, A system for crystallography and NMR, Yale University, New Haven CT) against 2.2Å data collected on an R-AXIS (Molecular Structure, Corp. Houston, TX) imaging plate detector.

f. Observations

As an effective recombinant human therapeutic, r-hu-G-CSF has been produced in large quantities and gram levels have been made available for structural analysis. The crystallization methods provided herein are likely to find other applications as other proteins of interest become available. This method can be applied to any crystallographic project which has large quantities of protein (approximately >200 mg). As one skilled in the art will recognize, the present materials and methods may be modified and equivalent materials and methods may be available for crystallization of other proteins.

B. Computer Program For Visualizing The Three Dimensional Structure of G-CSF

Although diagrams, such as those in the Figures herein, are useful for visualizing the three dimensional structure of G-CSF, a computer program which allows for stereoscopic viewing of the molecule is contemplated as preferred. This stereoscopic viewing, or "virtual reality" as those in the art sometimes refer to it, allows one to visualize the structure in its three dimensional form from every angle in a wide range of resolution, from macromolecular structure down to the atomic level. The computer programs contemplated herein also allow one to change perspective of the viewing angle of the molecule, for example by rotating the molecule. The contemplated programs also respond to changes so that one may, for example, delete, add, or substitute one or more images of atoms, including entire amino acid residues, or add chemical moieties to existing or substituted groups, and visualize the change in structure.

Other computer based systems may be used; the elements being: (a) a means for entering information, such as orthogonal coordinates or other numerically assigned coordinates of the three dimensional structure of G-CSF; (b) a means for expressing such coordinates, such as visual means so that one may view the three dimensional structure and correlate such three dimensional structure with the composition of the G-CSF molecule, such as the amino acid composition; (c) optionally, means for entering information which alters the composition of the G-CSF molecule expressed, so that the image of such three dimensional structure displays the altered composition.

The coordinates for the preferred computer program used are presented in FIGURE 5. The preferred computer program is Insight II, version 4, available from Biosym in San Diego, CA. For the raw crystallographic structure, the observed intensities of the diffraction data ("F-obs") and the orthogonal coordinates are also deposited in the Protein Data Bank, Chemistry Department, Brookhaven National Laboratory, Upton, New York 119723, USA and these are herein incorporated by reference.

Once the coordinates are entered into the Insight II program, one can easily display the three dimensional G-CSF molecule representation on a computer screen. The preferred computer system for display is Silicon Graphics 320 VGX (San Diego, CA). For stereoscopic viewing, one may wear eyewear (Crystal Eyes, Silicon Graphics) which allows one to visualize the G-CSF molecule in three dimensions stereoscopically, so one may turn the molecule and envision molecular design.

Thus, the present invention provides a method of designing or preparing a G-CSF analog with the aid of a computer comprising:

- (a) providing said computer with the means for displaying the three dimensional structure of a G-CSF molecule including displaying the composition of moieties of said G-CSF molecule, preferably displaying the three dimensional location of each amino acid, and more preferably displaying the three dimensional location of each atom of a G-CSF molecule;
- (b) viewing said display;
- (c) selecting a site on said display for alteration in the composition of said molecule or the location of a moiety; and
- (d) preparing a G-CSF analog with such alteration.

The alteration may be selected based on the desired structural characteristics of the end-product G-CSF analog, and considerations for such design are described in more detail below. Such considerations include the location and compositions of hydrophobic amino acid residues, particularly residues internal to the helical structures of a G-CSF molecule which residues, when altered, alter the overall structure of the internal core of the molecule and may prevent receptor binding; the location and compositions of external

loop structures, alteration of which may not affect the overall structure of the G-CSF molecule.

FIGURES 2-4 illustrate the overall three dimensional conformation in different ways. The topological diagram, the ribbon diagram, and the barrel diagram all illustrate aspects of the conformation of G-CSF.

FIGURE 2 illustrates a comparison between G-CSF and other molecules. There is a similarity of architecture, although these growth factors differ in the local conformations of their loops and bundle geometrics. The up-up-down-down topology with two long crossover connections is conserved, however, among all six of these molecules, despite the dissimilarity in amino acid sequence.

FIGURE 3 illustrates in more detail the secondary structure of recombinant human G-CSF. This ribbon diagram illustrates the handedness of the helices and their positions relative to each other.

FIGURE 4 illustrates in a different way the conformation of recombinant human G-CSF. This "barrel" diagram illustrates the overall architecture of recombinant human G-CSF.

C. Preparation of Analogs Using M13 Mutagenesis

This example relates to the preparation of G-CSF analogs using site directed mutagenesis techniques involving the single stranded bacteriophage M13, according to methods published in PCT Application No. WO 85/00817 (Souza et al., published February 28, 1985, herein incorporated by reference). This method essentially involves using a single-stranded nucleic acid template of the non-mutagenized sequence, and binding to it a smaller oligonucleotide containing the desired change in the sequence. Hybridization conditions allow for non-identical sequences to hybridize and the remaining sequence is filled in to be identical to the original template. What results is a double stranded molecule, with one of the two strands containing the desired change. This mutagenized single strand is separated, and used itself as a template for its complementary strand. This creates a double stranded molecule with the desired change.

The original G-CSF nucleic acid sequence used is presented in FIGURE 1, and the oligonucleotides containing the mutagenized nucleic acid(s) are presented in Table 2. Abbreviations used herein for amino acid residues and nucleotides are conventional, see Stryer, Biochemistry, 3d Ed., W.H. Freeman and Company, N.Y., N.Y. 1988, inside back cover.

The original G-CSF nucleic acid sequence was first placed into vector M13mp21. The DNA from single stranded phage M13mp21 containing the original G-CSF sequence was then isolated, and resuspended in water. For each reaction, 200 ng of this DNA was mixed with a 1.5 pmole of phosphorylated oligonucleotide (Table 2) and suspended in 0.1M Tris, 0.01M MgCl₂, 0.005M DTT, 0.1mM ATP, pH 8.0. The DNAs were annealed by heating to 65 °C and slowly cooling to room temperature.

Once cooled, 0.5mM of each ATP, dATP, dCTP, dGTP, TTP, 1 unit of T4 DNA ligase and 1 unit of Klenow fragment of *E. coli* polymerase 1 were added to the 1 unit of annealed DNA in 0.1M Tris, 0.025M NaCl, 0.01M MgCl₂, 0.01M DTT, pH 7.5.

The now double stranded, closed circular DNA was used to transfect *E. coli* without further purification. Plaques were screened by lifting the plaques with nitrocellulose filters, and then hybridizing the filters with single stranded DNA end-labeled with P³² for 1 hour at 55-60 °C. After hybridization, the filters were washed at 0-3 °C below the melt temperature of the oligo (2 °C for A-T, 4 °C for G-C) which selectively left autoradiography signals corresponding to plaques with phage containing the mutated sequence. Positive clones were confirmed by sequencing.

Set forth below are the oligonucleotides used for each G-CSF analog prepared via the M13 mutagenesis method. The nomenclature indicates the residue and the position of the original amino acid (e.g., Lysine at position 17), and the residue and position of the substituted amino acid (e.g., arginine 17). A substitution involving more than one residue is indicated via superscript notation, with commas between the noted positions or a semicolon indicating different residues. Deletions with no substitutions are so noted. The oligonucleotide sequences used for M13-based mutagenesis are next indicated; these oligonucleotides were manufactured synthetically, although the method of preparation is not critical, any nucleic acid synthesis method and/or equipment may be used. The length of the oligo is also indicated. As indicated above, these oligos were allowed to contact the single stranded phage vector, and then single nucleotides were added to complete the G-CSF analog nucleic acid sequence.

Table 2

G-CSE ANALOGS	SEQUENCES (5' -> 3')	Length (nucleotide)	Seq. ID
Lys17->Arg17	CTT TCT GCT GCG TTG TCT GGA ACA	24	3
Lys24->Arg24	ACA GGT TCG TCG TAT CCA GGG TG	23	4
Lys35->Arg35	CAC TGC AAG AAC GTC TGT GCG CT	23	5
Lys41->Arg41	CGC TAC TTA CCG TCT GTG CCA TC	23	6
Lys17, 24, 35-> Arg17, 24, 35	CTT TCT GCT GCG TTG TCT GGA ACA ACA GGT TCG TCG TAT CCA GGG TG CAC TGC AAG AAC GTC TGT GCG CT	24 23 23	7 8 9
Lys17, 24, 41-> Arg17, 24, 41	CTT TCT GCT GCG TTG TCT GGA ACA ACA GGT TCG TCG TAT CCA GGG TG CGC TAC TTA CCG TCT GTC CCA TC	24 23 23	10 11 12
Lys17, 35, 41-> Arg17, 35, 41	CTT TCT GCT GCG TTG TCT GGA ACA CAC TGC AAG AAC GTC TGT GCG CT CGC TAC TTA CCG TCT GTG CCA TC	24 23 23	13 14 15
Lys24, 35, 41-> Arg24, 35, 41	ACA GGT TCG TCG TAT CCA GGG TG CAC TGC AAG AAC GTC TGT GCG CT CGC TAC TTA CCG TCT GTG CCA TC	23 23 23	16 17 18

Table 2 (con't.)

G-CSE ANALOGS	SEQUENCES(5'→3')	Length(nucleotide)	Seq. ID
Lys ^{17,24,35,41} -> Arg ^{17,24,35,41}	CTT TCT GCT GCG TTG TCT GGA ACA ACA GGT TCG TCG TAT CCA GGG TG CAC TGC AAG AAC GTC TGT GCG CT CGC TAC TTA CCG TCT GTG CCA TC	24 23 23 23	19 20 21 22
Cys ¹⁸ ->Ala ¹⁸ Gln ⁶⁸ ->Glu ⁶⁸ Cys ^{37,43} -> Ser ^{37,43}	TCT GCT GAA AGC TCT GGA ACA GG CTT GTC CAT CTG AAG CTC TTC AG GAA AAA CTG TCC GCT ACT TAC AAA CTG TCC CAT CCG G	23 23 37	23 24 25
Gln ²⁶ ->Ala ²⁶ Gln ¹⁷⁴ ->Ala ¹⁷⁴	TTC GTA AAA TCG CCG GTG ACG G TCA TCT GGC TGC GCC GTA ATA G	22 22	26 27
Arg ¹⁷⁰ ->Ala ¹⁷⁰	CCG TGT TCT GGC TCA TCT GGC T	22	28
Arg ¹⁶⁷ ->Ala ¹⁶⁷	GAA GTA TCT TAC GCT GTT CTG CGT	24	29
Deletion 167	GAA GTA TCT TAC TAA GTT CTG CGT C	25	30
Lys ⁴¹ ->Ala ⁴¹	CGC TAC TTA CCG ACT GTG CCA T	22	31
His ⁴⁴ ->Lys ⁴⁴	CAA ACT GTG CAA GCC GGA AGA G	22	32
Glu ⁴⁷ ->Ala ⁴⁷	CAT CCG GAA GCA CTG GTA CTG C	22	33

Table 2 (con't.)

<u>G-CSE ANALOGS</u>	<u>SEQUENCES(5'--> 3')</u>	<u>Length(nucleotide)</u>	<u>Seq. ID</u>
Arg ²³ ->Ala ²³	GGA ACA GGT TGC TAA AAT CCA GG	23	34
Lys ²⁴ ->Ala ²⁴	GAA CAG GTT CGT GCG ATC CAG GGT G	25	35
Glu ²⁰ ->Ala ²⁰	GAA ATG TCT GGC ACA GGT TCG T	22	36
Asp ²⁸ ->Ala ²⁸	TCC AGG GTG CCG GTG CTG C	19	37
Met ¹²⁷ ->Glu ¹²⁷	AAG AGC TCG GTG AGG CAC CAG CT	23	38
Met ¹³⁸ ->Glu ¹³⁸	CTC AAG GTG CTG AGC CGG CAT TC	23	39
Met ¹²⁷ ->Leu ¹²⁷	GAG CTC GGT CTG GCA CCA GC	20	40
Met ¹³⁸ ->Leu ¹³⁸	TCA AGG TGC TCT GCC GGC ATT	21	41
Ser ¹³ ->Ala ¹³	TCT GCC GCA AGC CTT TCT GCT GA	23	42
Lys ¹⁷ ->Ala ¹⁷	CTT TCT GCT GGC ATG TCT GGA ACA	24	43
Gln ¹²¹ ->Ala ¹²¹	CTA TTT GGC AAG CGA TGG AAG AGC	24	44
Glu ¹²⁴ ->Ala ¹²⁴	CAG ATG GAA GCG CTC GGT ATG	21	45

Table 2 (con'tl)

G-CSF ANALOGS	SEQUENCES (5'→3')	Length(nucleotides)	Seq. ID
Met127,138→ Leu127,138	GAG CTC GGT CTG GCA CCA GC TCA AGG TGC TCT GCC GGC ATT	20 21	46 47
**Glu20→Ala20; Ser13→Gly13	GAA ATG TCT GGC ACA GGT TCG T	22	48

** This analog came about during the preparation of G-CSF analog Glu20→Ala20. As several clones were being sequenced to identify the Glu20→Ala20 analog, the Glu20→Ala20; Ser13→Gly13 analog was identified. This double mutant was the result of an in vitro Klenow DNA polymerase reaction mistake.

55 D. Preparation of G-CSF Analogs Using DNA Amplification

This example relates to methods for producing G-CSF analogs using a DNA amplification technique. Essentially, DNA encoding each analog was amplified in two separate pieces, combined, and then the total

sequence itself amplified. Depending upon where the desired change in the original G-CSF DNA was to be made, internal primers were used to incorporate the change, and generate the two separate amplified pieces. For example, for amplification of the 5' end of the desired analog DNA, a 5' flanking primer (complementary to a sequence of the plasmid upstream from the G-CSF original DNA) was used at one end of the region to be amplified, and an internal primer, capable of hybridizing to the original DNA but incorporating the desired change, was used for priming the other end. The resulting amplified region stretched from the 5' flanking primer through the internal primer. The same was done for the 3' terminus, using a 3' flanking primer (complementary to a sequence of the plasmid downstream from the G-CSF original DNA) and an internal primer complementary to the region of the intended mutation. Once the two "halves" (which may or may not be equal in size, depending on the location of the internal primer) were amplified, the two "halves" were allowed to connect. Once connected, the 5' flanking primer and the 3' flanking primer were used to amplify the entire sequence containing the desired change.

If more than one change is desired, the above process may be modified to incorporate the change into the internal primer, or the process may be repeated using a different internal primer. Alternatively, the gene amplification process may be used with other methods for creating changes in nucleic acid sequence, such as the phage based mutagenesis technique as described above. Examples of process for preparing analogs with more than one change are described below.

To create the G-CSF analogs described below, the template DNA used was the sequence as in FIGURE 1 plus certain flanking regions (from a plasmid containing the G-CSF coding region). These flanking regions were used as the 5' and 3' flanking primers and are set forth below. The amplification reactions were performed in 40 ul volumes containing 10 mM Tris-HCl, 1.5 mM MgCl₂, 50 mM KCl, 0.1 mg/ml gelatin, pH 8.3 at 20°C. The 40 ul reactions also contained 0.1mM of each dNTP, 10 pmoles of each primer, and 1 ng of template DNA. Each amplification was repeated for 15 cycles. Each cycle consisted of 0.5 minutes at 94°C, 0.5 minutes at 50°C, and 0.75 minutes at 72°C. Flanking primers were 20 nucleotides in length and internal primers were 20 to 25 nucleotides in length. This resulted in multiple copies of double stranded DNA encoding either the front portion or the back portion of the desired G-CSF analog.

For combining the two "halves," one fortieth of each of the two reactions was combined in a third DNA amplification reaction. The two portions were allowed to anneal at the internal primer location, as their ends bearing the mutation were complementary, and following a cycle of polymerization, give rise to a full length DNA sequence. Once so annealed, the whole analog was amplified using the 5' and 3' flanking primers. This amplification process was repeated for 15 cycles as described above.

The completed, amplified analog DNA sequence was cleaved with XbaI and XhoI restriction endonuclease to produce cohesive ends for insertion into a vector. The cleaved DNA was placed into a plasmid vector, and that vector was used to transform *E. coli*. Transformants were challenged with kanamycin at 50 ug/ml and incubated at 30°C. Production of G-CSF analog protein was confirmed by polyacrylamide gel electrophoresis of a whole cell lysate. The presence of the desired mutation was confirmed by DNA sequence analysis of plasmid purified from the production isolate. Cultures were then grown, and cells were harvested, and the G-CSF analogs were purified as set forth below.

Set forth below in Table 3 are the specific primers used for each analog made using gene amplification.

Table 3

Analog Seq. ID	Internal Primer(5'->3')	
His ⁴⁴ ->Ala ⁴⁴	5'primer-TTCCGGAGCGCACAGTTTG 3'primer-CAAACGTGGGGCTCCGGAAGAGC	49 50
Thr ¹¹⁷ ->Ala ¹¹⁷	5'primer-ATGCCAAATTGCAGTAGCAAAG 3'primer-CTTTGCTACTGCAATTTGGCAACA	51 52
Asp ¹¹⁰ ->Ala ¹¹⁰	5'primer-ATCAGCTACTGCTAGCTGCAGA 3'primer-TCTGCAGCTAGCAGTAGCTGACT	53 54
Gln ²¹ ->Ala ²¹	5'primer-TTACGAACCGCTTCCAGACATT 3'primer-AATGTCTGGAAGCGGTTTCGTAAAAT	55 56
Asp ¹¹³ ->Ala ¹¹³	5'primer-GTAGCAAATGCAGCTACATCTA 3'primer-TAGATGTAGCTGCATTTGCTACTAC	57 58
His ⁵³ ->Ala ⁵³	5'primer-CCAAGAGAAGCACCCAGCAG 3'primer-CTGCTGGGTGCTTCTCTTGGA	59 60
For each analog, the following 5' flanking primer was used:		
5'-CACTGGCGGTGATAATGAGC		61
For each analog, the following 3' flanking primer was used:		
3'-GGTCATTACGGACCGGATC		62

1. Construction of Double Mutation

To make G-CSF analog Gln^{12,21}->Glu^{12,21}, two separate DNA amplifications were conducted to create the two DNA mutations. The template DNA used was the sequence as in FIGURE 1 plus certain flanking regions (from a plasmid containing the G-CSF coding region). The precise sequences are listed below. Each of the two DNA amplification reactions were carried out using a Perkin Elmer/Cetus DNA Thermal Cycler. The 40 ul reaction mix consisted of 1X PCR Buffer (Cetus), 0.2 mM each of the 4 dXTPs (Cetus), 50 pmols of each primer oligonucleotide, 2 ng of G-CSF template DNA (on a plasmid vector), and 1 unit of Taq polymerase (Cetus). The amplification process was carried out for 30 cycles. Each cycle consisted of 1 minute at 94 °C, 2 minutes at 50 °C, and 3 minutes at 72 °C.

DNA amplification "A" used the oligonucleotides:

5' CCACTGGCGGTGATACTGAGC 3' (Seq. ID 63) and
5' AGCAGAAAGCTTTCCGGCAGAGAAGAAGCAGGA 3' (Seq. ID 64)
DNA amplification "B" used the oligonucleotides: 5' GCCGCAAAGCTTTCTGCTGAAATGTCTG-
GAAGAGGTTTCGTAAAATCCAGGGTGA 3' (Seq. ID 65) and
5' CTGGAATGCAGAAGCAAATGCCGGCATAGCACCTTCAGTCGGTTGCAGAGCTGGTGCCA 3' (Seq. ID 66)

From the 109 base pair double stranded DNA product obtained after DNA amplification "A", a 64 base pair XbaI to HindIII DNA fragment was cut and isolated that contained the DNA mutation Gln¹²->Glu¹². From the 509 base pair double stranded DNA product obtained after DNA amplification "B", a 197 base pair HindIII to BsmI DNA fragment was cut and isolated that contained the DNA mutation Gln²¹->Glu²¹.

The "A" and "B" fragments were ligated together with a 4.8 kilo-base pair XbaI to BsmI DNA plasmid vector fragment. The ligation mix consisted of equal molar DNA restriction fragments, ligation buffer (25 mM Tris-HCl pH 7.8, 10 mM MgCl₂, 2 mM DTT, 0.5 mM rATP, and 100 ug/ml BSA) and T4 DNA ligase and was incubated overnight at 14 °C. The ligated DNA was then transformed into *E. coli* FM5 cells by electroporation using a Bio Rad Gene Pulsar apparatus (BioRad, Richmond, CA). A clone was isolated and the plasmid construct verified to contain the two mutations by DNA sequencing. This 'intermediate' vector also contained a deletion of a 193 base pair BsmI to BsmI DNA fragment. The final plasmid vector was constructed by ligation and transformation (as described above) of DNA fragments obtained by cutting and isolating a 2 kilo-base pair SstI to BamHI DNA fragment from the intermediate vector, a 2.8 kbp SstI to EcoRI DNA fragment from the plasmid vector, and a 360 bp BamHI to EcoRI DNA fragment from the

plasmid vector. The final construct was verified by DNA sequencing the G-CSF gene. Cultures were grown, and the cells were harvested, and the G-CSF analogs were purified as set forth below.

As indicated above, any combination of mutagenesis techniques may be used to generate a G-CSF analog nucleic acid (and expression product) having one or more than one alteration. The two examples above, using M13-based mutagenesis and gene amplification-based mutagenesis, are illustrative.

E. Expression of G-CSF Analog DNA

The G-CSF analog DNAs were then placed into a plasmid vector and used to transform *E. coli* strain FM5 (ATCC#53911). The present G-CSF analog DNAs contained on plasmids and in bacterial host cells are available from the American Type Culture Collection, Rockville, MD, and the accession designations are indicated below.

One liter cultures were grown in broth containing 10g tryptone, 5g yeast extract and 5g NaCl) at 30 °C until reaching a density at A₆₀₀ of 0.5, at which point they were rapidly heated to 42 °C. The flasks were allowed to continue shaking at for three hours.

Other prokaryotic or eukaryotic host cells may also be used, such as other bacterial cells, strains or species, mammalian cells in culture (COS, CHO or other types) insect cells or multicellular organs or organisms, or plant cells or multicellular organs or organisms, and a skilled practitioner will recognize the appropriate host. The present G-CSF analogs and related compositions may also be prepared synthetically, as, for example, by solid phase peptide synthesis methods, or other chemical manufacturing techniques. Other cloning and expression systems will be apparent to those skilled in the art.

F. Purification of G-CSF Analog Protein

Cells were harvested by centrifugation (10,000 x G, 20 minutes, 4 °C). The pellet (usually 5 grams) was resuspended in 30 ml of 1mM DTT and passed three times through a French press cell at 10,000 psi. The broken cell suspension was centrifuged at 10,000g for 30 minutes, the supernatant removed, and the pellet resuspended in 30-40 ml water. This was recentrifuged at 10,000 x G for 30 minutes, and this pellet was dissolved in 25 ml of 2% Sarkosyl and 50mM Tris at pH 8. Copper sulfate was added to a concentration of 40uM, and the mixture was allowed to stir for at least 15 hours at 15-25 °C. The mixture was then centrifuged at 20,000 x G for 30 minutes. The resultant solubilized protein mixture was diluted four-fold with 13.3 mM Tris, pH 7.7, the Sarkosyl was removed, and the supernatant was then applied to a DEAE-cellulose (Whatman DE-52) column equilibrated in 20mM Tris, pH 7.7. After loading and washing the column with the same buffer, the analogs were eluted with 20mM Tris /NaCl (between 35mM to 100mM depending on the analog, as indicated below), pH 7.7. For most of the analogs, the eluent from the DEAE column was adjusted to a pH of 5.4, with 50% acetic acid and diluted as necessary (to obtain the proper conductivity) with 5mM sodium acetate pH 5.4. The solution was then loaded onto a CM-sepharose column equilibrated in 20 mM sodium acetate, pH 5.4. The column was then washed with 20mM NaAc, pH 5.4 until the absorbance at 280 nm was approximately zero. The G-CSF analog was then eluted with sodium acetate/NaCl in concentrations as described below in Table 4. The DEAE column eluents for those analogs not applied to the CM-sepharose column were dialyzed directly into 10mM NaAc, pH 4.0 buffer. The purified G-CSF analogs were then suitably isolated for *in vitro* analysis. The salt concentrations used for eluting the analogs varied, as noted above. Below, the salt concentrations for the DEAE cellulose column and for the CM-sepharose column are listed:

Table 4
Salt Concentrations

5	<u>Analog</u>	<u>DEAE Cellulose</u>	<u>CM-Sephadex</u>
	Lys17->Arg17	35mM	37.5mM
	Lys24->Arg24	35mM	37.5mM
10	Lys35->Arg35	35mM	37.5mM
	Lys41->Arg41	35mM	37.5mM
	Lys17, 24, 35-	35mM	37.5mM
15	>Arg17, 24, 35		
	Lys17, 35, 41-	35mM	37.5mM
	>Arg17, 35, 41		

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Table 4 Con't

	<u>Analog</u>	<u>DEAE Cellulose</u>	<u>CM-Sepharose</u>
5	Lys24, 35, 41-	35mM	37.5mM
	>Arg24, 35, 41		
10	Lys17, 24, 35, 41	35mM	37.5mM
	->Arg17, 24, 35, 41		
	Lys17, 24, 41-	35mM	37.5mM
	>Arg17, 24, 41		
15	Gln68->Glu68	60mM	37, 5mM
	Cys37, 43->Ser37, 43	40mM	37.5mM
	Gln26->Ala26	40mM	40mM
20	Gln174->Ala174	40mM	40mM
	Arg170->Ala170	40mM	40mM
	Arg167->Ala167	40mM	40mM
	Deletion 167*	N/A	N/A
25	Lys41->Ala41	160mM	40mM
	His44->Lys44	40mM	60mM
	Glu47->Ala47	40mM	40mM
30	Arg23->Ala23	40mM	40mM
	Lys24->Ala24	120mM	40mM
	Glu20->Ala20	40mM	60mM
35	Asp28->Ala28	40mM	80mM
	Met127->Glu127	80mM	40mM
	Met138->Glu138	80mM	40mM
	Met127->Leu127	40mM	40mM
40	Met138->Leu138	40mM	40mM
	Cys18->Ala18	40mM	37.5mM
	Gln12, 21->Glu12, 21	60mM	37.5mM
45	Gln12, 21, 68-	60mM	37.5mM
	>Glu12, 21, 68		
	Glu20->Ala20;		
	Ser13		
50	->Gly13	40mM	80mM

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Table 4 Con't

	<u>Analog</u>	<u>DEAE Cellulose</u>	<u>CM-Sephadex</u>
5	Met127,138-	40mM	40mM
	>Leu127,138		
10	Ser13->Ala13	40mM	40mM
	Lys17->Ala17	80mM	40mM
	Gln121->Ala121	40mM	60mM
15	Gln21->Ala21	50mM	Gradient 0 -150mM
	His44->Ala44**	40mM	N/A
	His53->Ala53**	50mM	N/A
	Asp110->Ala110**	40mM	N/A
20	Asp113->Ala113**	40mM	N/A
	Thr117->Ala117**	50mM	N/A
	Asp28->Ala28;	50mM	N/A
25	Asp110		
	Ala110**		
	Glu124->Ala124**	40mM	40mM

- 30 * For Deletion 167, the data are unavailable.
 ** For these analogs, the DEAE cellulose column alone
 was used for purification.

35 The above purification methods are illustrative, and a skilled practitioner will recognize that other means are available for obtaining the present G-CSF analogs.

G. Biological Assays

40 Regardless of which methods were used to create the present G-CSF analogs, the analogs were subject to assays for biological activity. Tritiated thymidine assays were conducted to ascertain the degree of cell division. Other biological assays, however, may be used to ascertain the desired activity. Biological assays such as assaying for the ability to induce terminal differentiation in mouse WEHI-3B (D+) leukemic cell line, also provides indication of G-CSF activity. See Nicola, et al., Blood 54: 614-27 (1979). Other in vitro assays may be used to ascertain biological activity. See Nicola, Annu. Rev. Biochem. 58: 45-77 (1989). In general, the test for biological activity should provide analysis for the desired result, such as increase or decrease in biological activity (as compared to non-altered G-CSF), different biological activity (as compared to non-altered G-CSF), receptor affinity analysis, or serum half-life analysis. The list is incomplete, and those skilled in the art will recognize other assays useful for testing for the desired end result.

50 The ³H-thymidine assay was performed using standard methods. Bone marrow was obtained from sacrificed female Balb C mice. Bone marrow cells were briefly suspended, centrifuged, and resuspended in a growth medium. A 160 ul aliquot containing approximately 10,000 cells was placed into each well of a 96 well micro-titer plate. Samples of the purified G-CSF analog(as prepared above) were added to each well, and incubated for 68 hours. Tritiated thymidine was added to the wells and allowed to incubate for 5 additional hours. After the 5 hour incubation time, the cells were harvested, filtered, and thoroughly rinsed. The filters were added to a vial containing scintillation fluid. The beta emissions were counted (LKB Betaplate scintillation counter). Standards and analogs were analyzed in triplicate, and samples which fell substantially above or below the standard curve were re-assayed with the proper dilution. The results

reported here are the average of the triplicate analog data relative to the unaltered recombinant human G-CSF standard results.

H. HPLC Analysis

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High pressure liquid chromatography was performed on purified samples of analog. Although peak position on a reverse phase HPLC column is not a definitive indication of structural similarity between two proteins, analogs which have similar retention times may have the same type of hydrophobic interactions with the HPLC column as the non-altered molecule. This is one indication of an overall similar structure.

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Samples of the analog and the non-altered recombinant human G-CSF were analyzed on a reverse phase (0.46 x 25 cm) Vydac 214TP54 column (Separations Group, Inc. Hesperia, CA). The purified analog G-CSF samples were prepared in 20 mM acetate and 40 mM NaCl solution buffered at pH 5.2 to a final concentration of 0.1 mg/ml to 5 mg/ml, depending on how the analog performed in the column. Varying amounts (depending on the concentration) were loaded onto the HPLC column, which had been equilibrated with an aqueous solution containing 1% isopropanol, 52.8% acetonitrile, and .38% trifluoro acetate (TFA).
15 The samples were subjected to a gradient of 0.86%/minute acetonitrile, and .002% TFA.

I. Results

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Presented below are the results of the above biological assays and HPLC analysis. Biological activity is the average of triplicate data and reported as a percentage of the control standard (non-altered G-CSF). Relative HPLC peak position is the position of the analog G-CSF relative to the control standard (non-altered G-CSF) peak. The "+" or "-" symbols indicate whether the analog HPLC peak was in advance of or followed the control standard peak (in minutes). Not all of the variants had been analyzed for relative HPLC
25 peak, and only those so analyzed are included below. Also presented are the American Type Culture Collection designations for E. coli host cells containing the nucleic acids coding for the present analogs, as prepared above.

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Table 5

Seq. ID	Variant	Analog	Relative		ATCC No.	% Normal G-CSF Activity
			HPLC Peak	Activity		
67	1	Lys17->Arg17	N/A	69184	N/A	
68	2	Lys24->Arg24	N/A	69185	N/A	
69	3	Lys35->Arg35	N/A	69186	N/A	
70	4	Lys41->Arg41	N/A	69187	N/A	
71	5	Lys17, 24, 35->Arg17, 24, 35	N/A	69189	N/A	
72	6	Lys17, 35, 41->Arg17, 35, 41	N/A	69192	N/A	
73	7	Lys24, 35, 41->Arg24, 35, 41	N/A	69191	N/A	
74	8	Lys17, 24, 35, 41 ->Arg17, 24, 35, 41	N/A	69193	N/A	
75	9	Lys17, 24, 41->Arg17, 24, 41	N/A	69190	N/A	
76	10	Gln68->Glu68	N/A	69196	N/A	
77	11	Cys37, 43->Ser37, 43	N/A	69197	N/A	
78	12	Gln26->Ala26	+ .96	69201	51%	
79	13	Gln174->Ala174	+ .14	69202	100%	
80	14	Arg170->Ala170	+ .78	69203	100%	

Table 5 Con't

Seq. ID	Variant	Analog	Relative		ATCC No.	% Normal
			HPLC Peak	G-CSF Activity		
81	15	Arg167->Ala167	+ .54	69204	110%	
82	16	Deletion 167	- .99	69207	N/A	
83	17	Lys41->Ala41	+ .25	69208	81%	
84	18	His44->Lys44	-1.53	69212	70%	
85	19	Glu47->Ala47	+ .14	69205	0%	
86	20	Arg23->Ala23	- .03	69206	31%	
87	21	Lys24->Ala24	+1.95	69213	0%	
88	22	Glu20->Ala20	-0.07	69211	0%	
89	23	Asp28->Ala28	- .30	69210	147%	
90	24	Met127->Glu127	N/A	69223	N/A	
91	25	Met138->Glu138	N/A	69222	N/A	
92	26	Met127->Leu127	N/A	69198	N/A	
93	27	Met138->Leu138	N/A	69199	N/A	
94	28	Cys18->Ala18	N/A	69188	N/A	
95	29	Gln12,21->Glu12,21	N/A	69194	N/A	
96	30	Gln12,21,68->Glu12,21,68	N/A	69195	N/A	
97	31	Glu20->Ala20; Ser13	+1.74	69209	0%	

Table 5 Con't

Seq. ID	Variant	Analog	Relative		ATCC No.	% Normal
			HPLC Peak	Activity		G-CSF
		->Gly13				
98	32	Met127,138->Leu127,138	+1.43	69200	98%	
99	33	Ser13->Ala13	0	69221	110%	
100	34	Lys17->Ala17	+0.50	69226	70%	
101	35	Gln121->Ala121	+2.7	69225	100%	
102	36	Gln21->Ala21	+0.63	69217	9.6%	
103	37	His44->Ala44	+1.52	69215	10.8%	
104	38	His53->Ala53	+0.99	69219	8.3%	
105	39	Asp110->Ala110	+1.97	69216	29%	
106	40	Asp113->Ala113	-0.34	69218	0%	
107	41	Thr117->Ala117	+0.4	69214	9.7%	
108	42	Asp28->Ala28; Asp110 Ala110	+3.2	69220	20.6%	

Table 5 Con't

Seq. ID	Variant	Analog	Relative		ATCC No.	% Normal	
			HPLC Peak	Activity		G-CSF	Activity
109	43	Glu124->Ala124	+0.16		69224		75%
110	44	Phe114->Val 114, Thr117->Ala117**	+0.53				0%

**This analog was apparently a result of an inadvertent error in the oligo which was used to prepare number 41, above (Thr117->Ala 117), and thus was prepared identically to the process used for that analog.

"N/A" indicates data which are not available.

1. Identification of Structure-Function Relationships

The first step used to design the present analogs was to determine what moieties are necessary for structural integrity of the G-CSF molecule. This was done at the amino acid residue level, although the

atomic level is also available for analysis. Modification of the residues necessary for structural integrity results in change in the overall structure of the G-CSF molecule. This may or may not be desirable, depending on the analog one wishes to produce. The working examples here were designed to maintain the overall structural integrity of the G-CSF molecule, for the purpose of maintain G-CSF receptor binding of the analog to the G-CSF receptor (as used in this section below, the "G-CSF receptor" refers to the natural G-CSF receptor, found on hematopoietic cells). It was assumed, and confirmed by the studies presented here, that G-CSF receptor binding is a necessary step for at least one biological activity, as determined by the above biological assays.

As can be seen from the figures, G-CSF (here, recombinant human met-G-CSF) is an antiparallel 4-alpha helical bundle with a left-handed twist, and with overall dimensions of 45 Å x 30Å x 24Å. The four helices within the bundle are referred to as helices A, B, C and D, and their connecting loops are known as the AB, BC and CD loops. The helix crossing angles range from -167.5° to -159.4°. Helices A, B, and C are straight, whereas helix D contains two kinds of structural characteristics, at Gly 150 and Ser 160 (of the recombinant human met-G-CSF). Overall, the G-CSF molecules is a bundle of four helices, connected in series by external loops. This structural information was then correlated with known functional information. It was known that residues (including methionine at position 1) 47, 23, 24, 20, 21, 44, 53, 113, 110, 28 and 114 may be modified, and the effect on biological activity would be substantial.

The majority of single mutations which lowered biological activity were centered around two regions of G-CSF that are separated by 30Å, and are located on different faces of the four helix bundle. One region involves interactions between the A helix and the D helix. This is further confirmed by the presence of salt bridges in the non-altered molecule as follows:

Atom	Helix	Atom	Helix	Distance
Arg 170 N1	D	Tyr 166 OH	A	3.3
Tyr 166 OH	D	Arg 23 N2	A	3.3
Glu 163 OE1	D	Arg 23 N1	A	2.8
Arg 23 N1	A	Gln 26 OE1	A	3.1
Gln 159 NE2	D	Gln 26 O	A	3.3

Distances reported here were for molecule A, as indicated in FIGURE 5 (wherein three G-CSF molecules crystallized together and were designated as A, B, and C). As can be seen, there is a web of salt bridges between helix A and helix D, which act to stabilize the helix A structure, and therefore affect the overall structure of the G-CSF molecule.

The area centering around residues Glu 20, Arg 23 and Lys 24 are found on the hydrophilic face of the A helix (residues 20-37). Substitution of the residues with the non-charged alanine residue at positions 20 and 23 resulted in similar HPLC retention times, indicating similarity in structure. Alteration of these sites altered the biological activity (as indicated by the present assays). Substitution at Lys 24 altered biological activity, but did not result in a similar HPLC retention time as the other two alterations.

The second site at which alteration lowered biological activity involves the AB helix. Changing glutamine at position 47 to alanine (analog no. 19, above) reduced biological activity (in the thymidine uptake assay) to zero. The AB helix is predominantly hydrophobic, except at the amino and carboxy termini; it contains one turn of a 3¹⁰ helix. There are two histadines at each termini (His 44 and His 56) and an additional glutamate at residue 46 which has the potential to form a salt bridge to His 44. The fourier transformed infra red spectrographic analysis (FTIR) of the analog suggests this analog is structurally similar to the non-altered recombinant G-CSF molecule. Further testing showed that this analog would not crystallize under the same conditions as the non-altered recombinant molecule.

Alterations at the carboxy terminus (Gln 174, Arg 167 and Arg 170) had little effect on biological activity. In contrast, deletion of the last eight residues (167-175) lowered biological activity. These results may indicate that the deletion destabilizes the overall structure which prevents the mutant from proper binding to the G-CSF receptor (and thus initiating signal transduction).

Generally, for the G-CSF internal core -- the internal four helix bundle lacking the external loops --the hydrophobic internal residues are essential for structural integrity. For example, in helix A, the internal hydrophobic residues are (with methionine being position 1) Phe 14, Cys 18, Val 22, Ile 25, Ile 32 and Leu 36. Generally, for the G-CSF internal core -- the internal four helix bundle lacking the external loops --the hydrophobic internal residues are essential for structural integrity. For example, in helix A, the internal hydrophobic residues are (with methionine being position 1 as in FIGURE 1) Phe 14, Cys 18, Val 22, Ile 25, Ile 32 and Leu 36. The other hydrophobic residues (again with the met at position 1) are: helix B, Ala 72,

Leu 76, Leu 79, Leu 83, Tyr 86, Leu 90, Leu 93; helix C, Leu 104, Leu 107, Val 111, Ala 114, Ile 118, Met 122; and helix D, Val 154, Val 158, Phe 161, Val 164, Val 168, Leu 172.

The above biological activity data, from the presently prepared G-CSF analogs, demonstrate that modification of the external loops interfere least with G-CSF overall structure. Preferred loops for analog preparation are the AB loop and the CD loop. The loops are relatively flexible structures as compared to the helices. The loops may contribute to the proteolysis of the molecule. G-CSF is relatively fast acting *in vivo* as the purpose the molecule serves is to generate a response to a biological challenge, i.e., selectively stimulate neutrophils. The G-CSF turnover rate is also relatively fast. The flexibility of the loops may provide a "handle" for proteases to attach to the molecule to inactivate the molecule. Modification of the loops to prevent protease degradation, yet have (via retention of the overall structure of non-modified G-CSF) no loss in biological activity may be accomplished.

This phenomenon is probably not limited to the G-CSF molecule but may also be common to the other molecules with known similar overall structures, as presented in Figure 2. Alteration of the external loop of, for example hGH, Interferon B, IL-2, GM-CSF and IL-4 may provide the least change to the overall structure. The external loops on the GM-CSF molecule are not as flexible as those found on the G-CSF molecule, and this may indicate a longer serum life, consistent with the broader biological activity of GM-CSF. Thus, the external loops of GM-CSF may be modified by releasing the external loops from the beta-sheet structure, which may make the loops more flexible (similar to those G-CSF) and therefore make the molecule more susceptible to protease degradation (and thus increase the turnover rate).

Alteration of these external loops may be effected by stabilizing the loops by connection to one or more of the internal helices. Connecting means are known to those in the art, such as the formation of a beta sheet, salt bridge, disulfide bonding or hydrophobic interactions, and other means are available. Also, deletion of one or more moieties, such as one or more amino acid residues or portions thereof, to prepare an abbreviated molecule and thus eliminate certain portions of the external loops may be effected.

Thus, by alteration of the external loops, preferably the AB loop (amino acids 58-72 of r-hu-met G-CSF) or the CD loop (amino acids 119 to 145 of r-hu-met-G-CSF), and less preferably the amino terminus (amino acids 1-10), one may therefore modify the biological function without elimination of G-CSF G-CSF receptor binding. For example, one may: (1) increase half-life (or prepare an oral dosage form, for example) of the G-CSF molecule by, for example, decreasing the ability of proteases to act on the G-CSF molecule or adding chemical modifications to the G-CSF molecule, such as one or more polyethylene glycol molecules or enteric coatings for oral formulation which would act to change some characteristic of the G-CSF molecule as described above, such as increasing serum or other half-life or decreasing antigenicity; (2) prepare a hybrid molecule, such as combining G-CSF with part or all of another protein such as another cytokine or another protein which effects signal transduction via entry through the cell through a G-CSF G-CSF receptor transport mechanism; or (3) increase the biological activity as in, for example, the ability to selectively stimulate neutrophils (as compared to a non-modified G-CSF molecule). This list is not limited to the above exemplars.

Another aspect observed from the above data is that stabilizing surface interactions may affect biological activity. This is apparent from comparing analogs 23 and 40. Analog 23 contains a substitution of the charged asparagine residue at position 28 for the neutrally-charged alanine residue in that position, and such substitution resulted in a 50% increase in the biological activity (as measured by the disclosed thymidine uptake assays). The asparagine residue at position 28 has a surface interaction with the asparagine residue at position 113; both residues being negatively charged, there is a certain amount of instability (due to the repelling of like charged moieties). When, however the asparagine at position 113 is replaced with the neutrally-charged alanine, the biological activity drops to zero (in the present assay system). This indicates that the asparagine at position 113 is critical to biological activity, and elimination of the asparagine at position 28 serves to increase the effect that asparagine at position 113 possesses.

The domains required for G-CSF receptor binding were also determined based on the above analogs prepared and the G-CSF structure. The G-CSF receptor binding domain is located at residues (with methionine being position 1) 11-57 (between the A and AB helix) and 100-118 (between the B and C helices). One may also prepare abbreviated molecules capable of binding to a G-CSF receptor and initiate signal transduction for selectively stimulating neutrophils by changing the external loop structure and having the receptor binding domains remain intact.

Residues essential for biological activity and presumably G-CSF receptor binding or signal transduction have been identified. Two distinct sites are located on two different regions of the secondary structure. What is here called "Site A" is located on a helix which is constrained by salt bridge contacts between two other members of the helical bundle. The second site, "Site B" is located on a relatively more flexible helix, AB. The AB helix is potentially more sensitive to local pH changes because of the type and position of the

residues at the carboxy and amino termini. The functional importance of this flexible helix may be important in a conformationally induced fit when binding to the G-CSF receptor. Additionally, the extended portion of the D helix is also indicated to be a G-CSF receptor binding domain, as ascertained by direct mutational and indirect comparative protein structure analysis. Deletion of the carboxy terminal end of r-hu-met-G-CSF reduces activity as it does for hGH, see, Cunningham and Wells, *Science* **244**: 1081-1084 (1989). Cytokines which have similar structures, such as IL-6 and GM-CSF with predicted similar topology also center their biological activity along the carboxy end of the D helix, see Bazan, *Immunology Today* **11**: 350-354 (1990).

A comparison of the structures and the positions of G-CSF receptor binding determinants between G-CSF and hGH suggests both molecules have similar means of signal transduction. Two separate G-CSF receptor binding sites have been identified for hGH De Vos et al., *Science* **255**: 306-32 (1991). One of these binding sites (called "Site I") is formed by residues on the exposed faces of hGH's helix 1, the connection region between helix 1 and 2, and helix 4. The second binding site (called "Site II") is formed by surface residues of helix 1 and helix 3.

The G-CSF receptor binding determinates identified for G-CSF are located in the same relative positions as those identified for hGH. The G-CSF receptor binding site located in the connecting region between helix A and B on the AB helix (Site A) is similar in position to that reported for a small piece of helix (residues 38-47) of hGH. A single point mutation in the AB helix of G-CSF significantly reduces biological activity (as ascertained in the present assays), indicating the role in a G-CSF receptor-ligand interface. Binding of the G-CSF receptor may destabilize the 3¹⁰ helical nature of this region and induce a conformation change improving the binding energy of the ligand/G-CSF receptor complex.

In the hGH receptor complex, the first helix of the bundle donates residues to both of the binding sites required to dimerize the hGH receptor. Mutational analysis of the corresponding helix of G-CSF (helix A) has identified three residues which are required for biological activity. Of these three residues, Glu 20 and Arg 24 lie on one face of the helical bundle towards helix C, whereas the side chain of Arg 23 (in two of the three molecules in the asymmetric unit) points to the face of the bundle towards helix D. The position of side chains of these biologically important residues indicates that similar to hGH, G-CSF may have a second G-CSF receptor binding site along the interface between helix A and helix C. In contrast with the hGH molecule, the amino terminus of G-CSF has a limited biological role as deletion of the first 11 residues has little effect on the biological activity.

As indicated above (see FIGURE 2, for example), G-CSF has a topological similarity with other cytokines. A correlation of the structure with previous biochemical studies, mutational analysis and direct comparison of specific residues of the hGH receptor complex indicates that G-CSF has two receptor binding sites. Site A lies along the interface of the A and D helices and includes residues in the small AB helix. Site B also includes residues in the A helix but lies along the interface between helices A and C. The conservation of structure and relative positions of biologically important residues between G-CSF and hGH is one indication of a common method of signal transduction in that the receptor is bound in two places. It is therefore found that G-CSF analogs possessing altered G-CSF receptor binding domains may be prepared by alteration at either of the G-CSF receptor binding sites (residues 20-57 and 145-175).

Knowledge of the three dimensional structure and correlation of the composition of G-CSF protein makes possible a systematic, rational method for preparing G-CSF analogs. The above working examples have demonstrated that the limitations of the size and polarity of the side chains within the core of the structure dictate how much change the molecule can tolerate before the overall structure is changed.

SEQUENCE LISTING

5 (1) GENERAL INFORMATION:

(i) APPLICANT: Amgen Inc.

(ii) TITLE OF INVENTION: G-CSF ANALOG COMPOSITIONS AND METHODS

10 (iii) NUMBER OF SEQUENCES: 110

(iv) CORRESPONDENCE ADDRESS:

(A) ADDRESSEE: Amgen Inc.

(B) STREET: Amgen Center, 1840 DeHavilland Drive

15 (C) CITY: Thousand Oaks

(D) STATE: California

(E) COUNTRY: United States of America

(F) ZIP: 91320-1789

(v) COMPUTER READABLE FORM:

20 (A) MEDIUM TYPE: Floppy disk

(B) COMPUTER: IBM PC compatible

(C) OPERATING SYSTEM: PC-DOS/MS-DOS

(2) INFORMATION FOR SEQ ID NO:1:

25 (i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 565 base pairs

(B) TYPE: nucleic acid

(C) STRANDEDNESS: single

(D) TOPOLOGY: linear

30 (ii) MOLECULE TYPE: DNA (genomic)

(ix) FEATURE:

(A) NAME/KEY: CDS

(B) LOCATION: 30..554

35 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:1:

TCTAGAAAAA	ACCAAGGAGG	TAATAAATA	ATG	ACT	CCA	TTA	GGT	CCT	GCT	TCT	53					
			Met	Thr	Pro	Leu	Gly	Pro	Ala	Ser						
			1				5									
40 TCT	CTG	CCG	CAA	AGC	TTT	CTG	CTG	AAA	TGT	CTG	GAA	CAG	GTT	CGT	AAA	101
Ser	Leu	Pro	Gln	Ser	Phe	Leu	Leu	Lys	Cys	Leu	Glu	Gln	Val	Arg	Lys	
	10					15				20						
45 ATC	CAG	GGT	GAC	GGT	GCT	GCA	CTG	CAA	GAA	AAA	CTG	TGC	GCT	ACT	TAC	149
Ile	Gln	Gly	Asp	Gly	Ala	Ala	Leu	Gln	Glu	Lys	Leu	Cys	Ala	Thr	Tyr	
25					30				35					40		

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EP 0 612 846 A1

	AAA CTG TGC CAT CCG GAA GAG CTG GTA CTG CTG GGT CAT TCT CTT GGG	197
	Lys Leu Cys His Pro Glu Glu Leu Val Leu Leu Gly His Ser Leu Gly	
	45 50 55	
5	ATC CCG TGG GCT CCG CTG TCT TCT TGT CCA TCT CAA GCT CTT CAG CTG	245
	Ile Pro Trp Ala Pro Leu Ser Ser Cys Pro Ser Gln Ala Leu Gln Leu	
	60 65 70	
	GCT GGT TGT CTG TCT CAA CTG CAT TCT GGT CTG TTC CTG TAT CAG GGT	293
10	Ala Gly Cys Leu Ser Gln Leu His Ser Gly Leu Phe Leu Tyr Gln Gly	
	75 80 85	
	CTT CTG CAA GCT CTG GAA GGT ATC TCT CCG GAA CTG GGT CCG ACT CTG	341
	Leu Leu Gln Ala Leu Glu Gly Ile Ser Pro Glu Leu Gly Pro Thr Leu	
	90 95 100	
15	GAC ACT CTG CAG CTA GAT GTA GCT GAC TTT GCT ACT ACT ATT TGG CAA	389
	Asp Thr Leu Gln Leu Asp Val Ala Asp Phe Ala Thr Thr Ile Trp Gln	
	105 110 115 120	
	CAG ATG GAA GAG CTC GGT ATG GCA CCA GCT CTG CAA CCG ACT CAA GGT	437
20	Gln Met Glu Glu Leu Gly Met Ala Pro Ala Leu Gln Pro Thr Gln Gly	
	125 130 135	
	GCT ATG CCG GCA TTC GCT TCT GCA TTC CAG CGT CGT GCA GGA GGT GTA	485
	Ala Met Pro Ala Phe Ala Ser Ala Phe Gln Arg Arg Ala Gly Gly Val	
	140 145 150	
25	CTG GTT GCT TCT CAT CTG CAA TCT TTC CTG GAA GTA TCT TAC CGT GTT	533
	Leu Val Ala Ser His Leu Gln Ser Phe Leu Glu Val Ser Tyr Arg Val	
	155 160 165	
	CTG CGT CAT CTG GCT CAG CCG TAATAGAATT C	565
30	Leu Arg His Leu Ala Gln Pro	
	170 175	

(2) INFORMATION FOR SEQ ID NO:2:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:2:

40	Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
	1 5 10 15
	Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
	20 25 30
45	Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
	35 40 45

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EP 0 612 846 A1

Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:3:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 24 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:3:

CTTTCTGCTG CGTGTCTGG AACA

24

(2) INFORMATION FOR SEQ ID NO:4:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 23 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:4:

ACAGGTTTGGT CGTATCCAGG GTG

23

(2) INFORMATION FOR SEQ ID NO:5:

- (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:5:

CACTGCAAGA ACGTCTGTGC GCT

23

(2) INFORMATION FOR SEQ ID NO:6:

- (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:6:

CGCTACTTAC CGTCTGTGCC ATC

23

(2) INFORMATION FOR SEQ ID NO:7:

- (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 24 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:7:

CTTTCTGCTG CGTTGTCTGG AACA

24

(2) INFORMATION FOR SEQ ID NO:8:

- (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:8:
ACAGGTTTCGT CGTATCCAGG GTG 23

5 (2) INFORMATION FOR SEQ ID NO:9:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 23 base pairs
(B) TYPE: nucleic acid
10 (C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:9:
15 CACTGCAAGA ACGTCTGTGC GCT 23

(2) INFORMATION FOR SEQ ID NO:10:

(i) SEQUENCE CHARACTERISTICS:
20 (A) LENGTH: 24 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

25 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:10:
CTTTCTGCTG CGTTGTCTGG AACA 24

30 (2) INFORMATION FOR SEQ ID NO:11:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 23 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
35 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:11:
40 ACAGGTTTCGT CGTATCCAGG GTG 23

(2) INFORMATION FOR SEQ ID NO:12:

(i) SEQUENCE CHARACTERISTICS:
45 (A) LENGTH: 23 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single

50

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(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:12:

5 CGCTACTTAC CGTCTGTCCC ATC 23

(2) INFORMATION FOR SEQ ID NO:13:

10 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 24 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

15 (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:13:

CTTTCTGCTG CGTTGTCTGG AACAA 24

20 (2) INFORMATION FOR SEQ ID NO:14:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 25 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:14:

30 CACTGCAAGA ACGTCTGTGC GCT 23

(2) INFORMATION FOR SEQ ID NO:15:

35 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

40 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:15:

CGCTACTTAC CGTCTGTGCC ATC 23

45

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(2) INFORMATION FOR SEQ ID NO:16:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 23 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:16:

ACAGGTTTCGT CGTATCCAGG GTG

23

(2) INFORMATION FOR SEQ ID NO:17:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 23 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:17:

CACTGCAAGA ACGTCTGTGC GCT

23

(2) INFORMATION FOR SEQ ID NO:18:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 23 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:18:

CGCTACTTAC CGTCTGTGCC ATC

23

(2) INFORMATION FOR SEQ ID NO:19:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 24 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:19:
 CTTTCTGCTG CGTTGTCTGG AACA 24

5 (2) INFORMATION FOR SEQ ID NO:20:
 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 10 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear
 (ii) MOLECULE TYPE: DNA
 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:20:
 15 ACAGGTTTCGT CGTATCCAGG GTG 23

(2) INFORMATION FOR SEQ ID NO:21:
 (i) SEQUENCE CHARACTERISTICS:
 20 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear
 (ii) MOLECULE TYPE: DNA
 25 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:21:
 CACTGCAAGA ACGTCTGTGC GCT 23

30 (2) INFORMATION FOR SEQ ID NO:22:
 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 35 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear
 (ii) MOLECULE TYPE: DNA
 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:22:
 40 CGCTACTTAC CGTCTGTGCC ATC 23

(2) INFORMATION FOR SEQ ID NO:23:
 (i) SEQUENCE CHARACTERISTICS:
 45 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single

50

55

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

5 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:23:

TCTGCTGAAA GCTCTGGAAC AGG 23

(2) INFORMATION FOR SEQ ID NO:24:

10 (i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 23 base pairs

(B) TYPE: nucleic acid

(C) STRANDEDNESS: single

(D) TOPOLOGY: linear

15 (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:24:

CTTGTCCATC TGAAGCTCTT CAG 23

20 (2) INFORMATION FOR SEQ ID NO:25:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 37 base pairs

(B) TYPE: nucleic acid

25 (C) STRANDEDNESS: single

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:25:

30 GAAAAACTGT CCGCTACTTA CAAACTGTCC CATCCGG 37

(2) INFORMATION FOR SEQ ID NO:26:

(i) SEQUENCE CHARACTERISTICS:

35 (A) LENGTH: 22 base pairs

(B) TYPE: nucleic acid

(C) STRANDEDNESS: single

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

40 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:26:

TTCGTAAAAT CGCGGGTGAC GG 22

45

50

55

(2) INFORMATION FOR SEQ ID NO:27:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 22 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:27:

TCATCTGGCT GCGCCGTAAT AG

22

(2) INFORMATION FOR SEQ ID NO:28:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 22 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:28:

CCGTGTTCTG GCTCATCTGG CT

22

(2) INFORMATION FOR SEQ ID NO:29:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 24 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:29:

GAAGTATCTT ACGCTGTTCT GCGT

24

(2) INFORMATION FOR SEQ ID NO:30:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 25 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:30:
GAAGTATCTT ACTAAGTTCT GCGTC 25

5 (2) INFORMATION FOR SEQ ID NO:31:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 22 base pairs
(B) TYPE: nucleic acid
10 (C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:31:
15 CGCTACTTAC GCACTGTGCC AT 22

(2) INFORMATION FOR SEQ ID NO:32:

(i) SEQUENCE CHARACTERISTICS:
20 (A) LENGTH: 22 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

25 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:32:
CAAACGTGTC AAGCCGGAAG AG 22

30 (2) INFORMATION FOR SEQ ID NO:33:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 22 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
35 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:33:
40 CATCCGGAAG CACTGGTACT GC 22

(2) INFORMATION FOR SEQ ID NO:34:

(i) SEQUENCE CHARACTERISTICS:
45 (A) LENGTH: 23 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single

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(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

5 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:34:

GGAACAGGTT GCTAAAATCC AGG 23

(2) INFORMATION FOR SEQ ID NO:35:

10 (i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 25 base pairs

(B) TYPE: nucleic acid

(C) STRANDEDNESS: single

(D) TOPOLOGY: linear

15 (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:35:

GAACAGGTTC GTGCGATCCA GGGTG 25

(2) INFORMATION FOR SEQ ID NO:36:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 22 base pairs

(B) TYPE: nucleic acid

25 (C) STRANDEDNESS: single

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:36:

30 GAAATGTCTG GCACAGGTTC GT 22

(2) INFORMATION FOR SEQ ID NO:37:

(i) SEQUENCE CHARACTERISTICS:

35 (A) LENGTH: 19 base pairs

(B) TYPE: nucleic acid

(C) STRANDEDNESS: single

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

40 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:37:

TCCAGGGTGC CGGTGCTGC 19

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(2) INFORMATION FOR SEQ ID NO:38:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 23 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:38:

AAGAGCTCGG TGAGGCACCA GCT

23

(2) INFORMATION FOR SEQ ID NO:39:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 23 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:39:

CTCAAGGTGC TGAGCCGGCA TTC

23

(2) INFORMATION FOR SEQ ID NO:40:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 20 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:40:

GAGCTCGGTC TGGCACCAGC

20

(2) INFORMATION FOR SEQ ID NO:41:

- (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 21 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:41:
TCAAGGTGCT CTGCCGGCAT T 21

5 (2) INFORMATION FOR SEQ ID NO:42:
 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
10 (D) TOPOLOGY: linear
 (ii) MOLECULE TYPE: DNA
 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:42:
15 TCTGCCGCAA GCCTTTCTGC TGA 23

(2) INFORMATION FOR SEQ ID NO:43:
 (i) SEQUENCE CHARACTERISTICS:
20 (A) LENGTH: 24 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear
 (ii) MOLECULE TYPE: DNA
25 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:43:
CTTTCTGCTG GCATGTCTGG AACA 24

30 (2) INFORMATION FOR SEQ ID NO:44:
 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 24 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
35 (D) TOPOLOGY: linear
 (ii) MOLECULE TYPE: DNA
 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:44:
40 CTATTTGGCA AGCGATGGAA GAGC 24

(2) INFORMATION FOR SEQ ID NO:45:
 (i) SEQUENCE CHARACTERISTICS:
45 (A) LENGTH: 21 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single

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(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

5 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:45:
CAGATGGAAG CGCTCGGTAT G 21

(2) INFORMATION FOR SEQ ID NO:46:

10 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 20 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

15 (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:46:
GAGCTCGGTC TGGCACCAGC 20

(2) INFORMATION FOR SEQ ID NO:47:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 21 base pairs
 (B) TYPE: nucleic acid
 25 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:47:
30 TCAAGGTGCT CTGCCGGCAT T 21

(2) INFORMATION FOR SEQ ID NO:48:

35 (i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 22 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

40 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:48:
GAAATGTCTG GCACAGGTTC GT 22

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(2) INFORMATION FOR SEQ ID NO:49:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 19 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:49:

TTCCGGAGCG CACAGTTTG

19

(2) INFORMATION FOR SEQ ID NO:50:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:50:

CGAGAAGGCC TCGGGTGTCA AAC

23

(2) INFORMATION FOR SEQ ID NO:51:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 22 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:51:

ATGCCAAATT GCAGTAGCAA AG

22

(2) INFORMATION FOR SEQ ID NO:52:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 24 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:52:
 5 ACAACGGTTT AACGTCATCG TTTC 24

(2) INFORMATION FOR SEQ ID NO:53:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 22 base pairs
 10 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:53:
 15 ATCAGCTACT GCTAGCTGCA GA 22

(2) INFORMATION FOR SEQ ID NO:54:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 23 base pairs
 20 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA
 25

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:54:
 TCAGTCGATG ACGATCGACG TCT 23

(2) INFORMATION FOR SEQ ID NO:55:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 22 base pairs
 30 (B) TYPE: nucleic acid
 35 (C) STRANDEDNESS: single
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:55:
 40 TTACGAACCG CTCCAGACA TT 22

(2) INFORMATION FOR SEQ ID NO:56:

(i) SEQUENCE CHARACTERISTICS:
 45 (A) LENGTH: 25 base pairs
 (B) TYPE: nucleic acid
 (C) STRANDEDNESS: single

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(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

5 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:56:
TAAAATGCTT GGCGAAGGTC TGTA 25

(2) INFORMATION FOR SEQ ID NO:57:

10 (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 22 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

15 (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:57:
GTAGCAAATG CAGCTACATC TA 22

20 (2) INFORMATION FOR SEQ ID NO:58:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 25 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
25 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:58:
30 CATCATCGTT TACGTCGATG TAGAT 25

(2) INFORMATION FOR SEQ ID NO:59:

(i) SEQUENCE CHARACTERISTICS:
35 (A) LENGTH: 20 base pairs
(B) TYPE: nucleic acid
(C) STRANDEDNESS: single
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

40 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:59:
CCAAGAGAAG CACCCAGCAG 20

45

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55

(2) INFORMATION FOR SEQ ID NO:60:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 22 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:60:

AGGGTTCTCT TCGTGGGTCG TC

22

(2) INFORMATION FOR SEQ ID NO:61:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 20 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:61:

CACTGGCGGT GATAATGAGC

20

(2) INFORMATION FOR SEQ ID NO:62:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 19 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:62:

CTAGGCCAGG CATTACTGG

19

(2) INFORMATION FOR SEQ ID NO:63:

- (i) SEQUENCE CHARACTERISTICS:
 - (A) LENGTH: 21 base pairs
 - (B) TYPE: nucleic acid
 - (C) STRANDEDNESS: single
 - (D) TOPOLOGY: linear

- (ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:63:

CCACTGGCGG TGATACTGAG C

21

5

(2) INFORMATION FOR SEQ ID NO:64:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 33 base pairs

(B) TYPE: nucleic acid

10 (C) STRANDEDNESS: single

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:64:

15

AGCAGAAAGC TTTCCGGCAG AGAAGAAGCA GGA

33

(2) INFORMATION FOR SEQ ID NO:65:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 54 base pairs

(B) TYPE: nucleic acid

20 (C) STRANDEDNESS: single

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

25

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:65:

GCCGCAAAGC TTTCTGCTGA AATGTCTGGA AGAGGTTCGT AAAATCCAGG GTGA

54

30

(2) INFORMATION FOR SEQ ID NO:66:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 59 base pairs

(B) TYPE: nucleic acid

35 (C) STRANDEDNESS: single

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: DNA

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:66:

40

CTGGAATGCA GAAGCAAATG CCGGCATAGC ACCTTCAGTC GGTTCAGAG CTGGTGCCA

59

(2) INFORMATION FOR SEQ ID NO:67:

(i) SEQUENCE CHARACTERISTICS:

45

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

(D) TOPOLOGY: linear

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(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:67:

5 Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
10 Arg Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
15 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
20 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
25 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
30 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:68:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:68:

45 Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
50 Lys Cys Leu Glu Gln Val Arg Arg Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30

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Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
5 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
10 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
15 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
20 Phe Gln Arg Arg Ala Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

25

(2) INFORMATION FOR SEQ ID NO:69:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

(D) TOPOLOGY: linear

30

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:69:

35

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15

Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30

40

Gln Glu Arg Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45

Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

45

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

50

55

EP 0 612 846 A1

Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
5 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
10 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

15

(2) INFORMATION FOR SEQ ID NO:70:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
20 (B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:70:

25

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
30 Gln Glu Lys Leu Cys Ala Thr Tyr Arg Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
35 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
40 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
45 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
50 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

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EP 0 612 846 A1

Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

5

(2) INFORMATION FOR SEQ ID NO:71:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

10

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:71:

15 Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Arg Cys Leu Glu Gln Val Arg Arg Ile Gln Gly Asp Gly Ala Ala Leu
20 20 25 30
Gln Glu Arg Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
25 35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
25 65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
30 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
35 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
40 145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

45 (2) INFORMATION FOR SEQ ID NO:72:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

50

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

55

• (xi) SEQUENCE DESCRIPTION: SEQ ID NO:72:

5 Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Arg Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
10 Gln Glu Arg Leu Cys Ala Thr Tyr Arg Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
15 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
20 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
25 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
30 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:73:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:73:

45 Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Arg Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
50 Gln Glu Arg Leu Cys Ala Thr Tyr Arg Leu Cys His Pro Glu Glu Leu
35 40 45

55

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Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

5 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

10 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110

Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125

15 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140

Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

20 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:74:

25 (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

30 (ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:74:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15

35 Arg Cys Leu Glu Gln Val Arg Arg Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30

Gln Glu Arg Leu Cys Ala Thr Tyr Arg Leu Cys His Pro Glu Glu Leu
35 40 45

40 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

45 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

50 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110

55

EP 0 612 846 A1

Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
5 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
10 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:75:

15 (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

20 (ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:75:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
25 Arg Cys Leu Glu Gln Val Arg Arg Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Arg Leu Cys His Pro Glu Glu Leu
30 35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
35 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
40 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
45 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
50 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

55

(2) INFORMATION FOR SEQ ID NO:76:

(i) SEQUENCE CHARACTERISTICS:

- 5 (A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

10

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:76:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 15 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 20 50 55 60
 Cys Pro Ser Glu Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 25 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 30 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 35 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 40 165 170 175

(2) INFORMATION FOR SEQ ID NO:77:

(i) SEQUENCE CHARACTERISTICS:

- 45 (A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

50

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:77:

55

EP 0 612 846 A1

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
5 20 25 30
Gln Glu Lys Leu Ser Ala Thr Tyr Lys Leu Ser His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
10 50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
15 85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
20 115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
25 145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:78:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
- (B) TYPE: amino acid
- (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:78:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Ala Gly Asp Gly Ala Ala Leu
45 20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 50 55 60

EP 0 612 846 A1

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 5 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 10 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 15 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

20 (2) INFORMATION FOR SEQ ID NO:79:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 25 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:79:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125

55

EP 0 612 846 A1

Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140

5 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Ala Pro
165 170 175

10

(2) INFORMATION FOR SEQ ID NO:80:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
15 (B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

20 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:80:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15

Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
25 20 25 30

Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45

Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

35 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110

40 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125

Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140

45 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

Phe Leu Glu Val Ser Tyr Arg Val Leu Ala His Leu Ala Gln Pro
50 165 170 175

55

(2) INFORMATION FOR SEQ ID NO:81:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:81:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Ala Val Leu Arg His Leu Ala Gln Pro
 165 170 175

(2) INFORMATION FOR SEQ ID NO:82:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 174 amino acids

(B) TYPE: amino acid

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:82:

EP 0 612 846 A1

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 5 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 10 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 15 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 20 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 25 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Val Leu Arg His Leu Ala Gln Pro
 165 170 174

30

(2) INFORMATION FOR SEQ ID NO:83:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 (D) TOPOLOGY: linear

35

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:83:

40

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 45 Gln Glu Lys Leu Cys Ala Thr Tyr Ala Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60

50

55

EP 0 612 846 A1

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 5 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 10 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 15 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175
 20

(2) INFORMATION FOR SEQ ID NO:84:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 175 amino acids
 25 (B) TYPE: amino acid
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

30 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:84:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys Lys Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 40 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 45 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 50 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 55

EP 0 612 846 A1

Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

(2) INFORMATION FOR SEQ ID NO:85:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:85:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Ala Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

EP 0 612 846 A1

(2) INFORMATION FOR SEQ ID NO:86:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:86:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Ala Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:87:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:87:

EP 0 612 846 A1

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
5 Lys Cys Leu Glu Gln Val Arg Ala Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
10 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
15 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
20 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
25 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175
30

(2) INFORMATION FOR SEQ ID NO:88:

(i) SEQUENCE CHARACTERISTICS:
35 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

40 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:88:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
45 Lys Cys Leu Ala Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
50 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

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EP 0 612 846 A1

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 5 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 10 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 15 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

(2) INFORMATION FOR SEQ ID NO:89:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:89:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Ala Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 45 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 50 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125

EP 0 612 846 A1

Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

(2) INFORMATION FOR SEQ ID NO:90:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
- (B) TYPE: amino acid
- (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:90:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Glu Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

EP 0 612 846 A1

(2) INFORMATION FOR SEQ ID NO:91:

(i) SEQUENCE CHARACTERISTICS:

- 5 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

10 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:91:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
15 20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Glu Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:92:

(i) SEQUENCE CHARACTERISTICS:

- 45 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

50 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:92:

55

EP 0 612 846 A1

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe L u Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Leu Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

30

(2) INFORMATION FOR SEQ ID NO:93:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

35

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:93:

40

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

50

55

EP 0 612 846 A1

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 5 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 10 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Leu Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 15 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

20

(2) INFORMATION FOR SEQ ID NO:94:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

25

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:94:

30

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15

Lys Ala Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30

35

Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45

Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60

40

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80

Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95

45

Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110

Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125

50

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Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

(2) INFORMATION FOR SEQ ID NO:95:

(i) SEQUENCE CHARACTERISTICS:

(A) LENGTH: 175 amino acids

(B) TYPE: amino acid

(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:95:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Glu Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Glu Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

(2) INFORMATION FOR SEQ ID NO:96:

(i) SEQUENCE CHARACTERISTICS:

- 5 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

10 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:96:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Glu Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Glu Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
15 20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
20 50 55 60
Cys Pro Ser Glu Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
25 85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
30 115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
35 130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
40 165 170 175

(2) INFORMATION FOR SEQ ID NO:97:

(i) SEQUENCE CHARACTERISTICS:

- 45 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

50 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:97:

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Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Gly Phe Leu Leu
 1      5      10      15
5  Lys Cys Leu Ala Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
      20      25      30
    Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
      35      40      45
10 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
   50      55      60
15 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
   65      70      75      80
    Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
      85      90      95
20 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
   100      105      110
    Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
      115      120      125
25 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
   130      135      140
    Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
   145      150      155      160
30 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
      165      170      175

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(2) INFORMATION FOR SEQ ID NO:98:

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35      (i) SEQUENCE CHARACTERISTICS:
          (A) LENGTH: 175 amino acids
          (B) TYPE: amino acid
          (D) TOPOLOGY: linear

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40      (ii) MOLECULE TYPE: protein

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(xi) SEQUENCE DESCRIPTION: SEQ ID NO:98:

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Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1      5      10      15
45 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
      20      25      30
50 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
      35      40      45

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Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

5 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

10 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110

Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Leu Ala
115 120 125

15 Pro Ala Leu Gln Pro Thr Gln Gly Ala Leu Pro Ala Phe Ala Ser Ala
130 135 140

Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

20 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:99:

25 (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

30 (ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:99:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ala Phe Leu Leu
1 5 10 15

35 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30

Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45

40 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

45 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

50 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110

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Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
5 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
10 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:100:

15 (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

20 (ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:100:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
25 Ala Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
30 35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
35 65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
40 100 105 110
Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
45 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
50 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

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(2) INFORMATION FOR SEQ ID NO:101:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
- (B) TYPE: amino acid
- (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:101:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Ala Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

(2) INFORMATION FOR SEQ ID NO:102:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
- (B) TYPE: amino acid
- (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:102:

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Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
5 Lys Cys Leu Glu Ala Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
10 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
15 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
20 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
25 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175
30

(2) INFORMATION FOR SEQ ID NO:103:

(i) SEQUENCE CHARACTERISTICS:
35 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

40 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:103:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
45 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys Ala Pro Glu Glu Leu
35 40 45
50 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

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Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 5 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 10 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 15 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175

20 (2) INFORMATION FOR SEQ ID NO:104:

(i) SEQUENCE CHARACTERISTICS:
 (A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 25 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:104:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 Val Leu Leu Gly Ala Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 40 65 70 75 80
 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 45 100 105 110
 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 50 115 120 125

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Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140

5 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

10

• (2) INFORMATION FOR SEQ ID NO:105:

(i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
15 (B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

20 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:105:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15

Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
25 20 25 30

Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45

30 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

35 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Ala Val Ala
100 105 110

40 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125

Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140

45 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

50 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

55

(2) INFORMATION FOR SEQ ID NO:106:

(i) SEQUENCE CHARACTERISTICS:

- 5 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

10 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:106:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Ala Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:107:

(i) SEQUENCE CHARACTERISTICS:

- 45 (A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

50 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:107:

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Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
5 20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
10 50 55 60
Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80
Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
15 85 90 95
Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110
Asp Phe Ala Thr Ala Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
20 115 120 125
Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140
Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
25 145 150 155 160
Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:108:

(i) SEQUENCE CHARACTERISTICS:

- (A) LENGTH: 175 amino acids
- (B) TYPE: amino acid
- (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:108:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15
Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Ala Gly Ala Ala Leu
45 20 25 30
Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45
Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

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Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 5 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Ala Val Ala
 100 105 110
 10 Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
 115 120 125
 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
 130 135 140
 15 Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
 145 150 155 160
 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
 165 170 175
 20

(2) INFORMATION FOR SEQ ID NO:109:

(i) SEQUENCE CHARACTERISTICS:
 25 (A) LENGTH: 175 amino acids
 (B) TYPE: amino acid
 (D) TOPOLOGY: linear

(ii) MOLECULE TYPE: protein

30 (xi) SEQUENCE DESCRIPTION: SEQ ID NO:109:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
 1 5 10 15
 35 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
 20 25 30
 Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
 35 40 45
 40 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
 50 55 60
 Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
 65 70 75 80
 45 Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
 85 90 95
 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
 100 105 110
 50

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Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Ala Leu Gly Met Ala
115 120 125

5 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140

Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

10 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

(2) INFORMATION FOR SEQ ID NO:110:

15 (i) SEQUENCE CHARACTERISTICS:
(A) LENGTH: 175 amino acids
(B) TYPE: amino acid
(D) TOPOLOGY: linear

20 (ii) MOLECULE TYPE: protein

(xi) SEQUENCE DESCRIPTION: SEQ ID NO:110:

Met Thr Pro Leu Gly Pro Ala Ser Ser Leu Pro Gln Ser Phe Leu Leu
1 5 10 15

25 Lys Cys Leu Glu Gln Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu
20 25 30

Gln Glu Lys Leu Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu
35 40 45

30 Val Leu Leu Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser
50 55 60

Cys Pro Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His
65 70 75 80

Ser Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
85 90 95

40 Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val Ala
100 105 110

Asp Val Ala Thr Ala Ile Trp Gln Gln Met Glu Glu Leu Gly Met Ala
115 120 125

45 Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe Ala Ser Ala
130 135 140

Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser His Leu Gln Ser
145 150 155 160

50 Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His Leu Ala Gln Pro
165 170 175

55 Claims

1. A method for preparing a G-CSF analog comprising the steps of:
 - (a) viewing information conveying the three dimensional structure of a G-CSF molecule;

- (b) selecting from said viewed information at least one site on said G-CSF molecule for alteration;
 - (c) preparing a G-CSF molecule having such alteration; and
 - (d) optionally, testing such G-CSF molecule for a desired characteristic.
- 5 2. A computer based method for preparing a G-CSF analog comprising the steps of:
- (a) providing computer expression of the three dimensional structure of a G-CSF molecule;
 - (b) selecting from said computer expression at least one site on said G-CSF molecule for alteration;
 - (c) preparing a G-CSF molecule having such alteration; and,
 - (d) optionally, testing such G-CSF molecule for a desired characteristic.
- 10 3. A method for preparing a G-CSF analog with the aid of a computer comprising:
- (a) providing said computer with the means for displaying the three dimensional structure of a G-CSF molecule including displaying the composition of moieties of said G-CSF molecule, preferably displaying the three dimensional location of each amino acid, and more preferably displaying the three dimensional location of each atom of a G-CSF molecule;
 - (b) viewing said display;
 - (c) selecting a site on said display for alteration in the composition of said molecule or the location of a moiety; and
 - (d) preparing a G-CSF analog with such alteration.
- 15 20 4. A computer-based method for preparing a G-CSF analog comprising the steps of:
- (a) viewing the three dimensional structure of a G-CSF molecule via a computer, said computer having been previously programmed (i) to express the coordinates of a G-CSF molecule in three dimensional space, and (ii) to allow for entry of information for alteration of said G-CSF expression and viewing thereof;
 - (b) selecting a site on said visual image of said G-CSF molecule for alteration;
 - (c) entering information for said alteration on said computer;
 - (d) viewing a three dimensional structure of said altered G-CSF molecule via said computer;
 - (e) optionally repeating steps (a)-(e) above;
 - (f) preparing a G-CSF analog with said alteration; and
 - (g) optionally testing said G-CSF analog for a desired characteristic.
- 25 30 5. In a computer-based apparatus for displaying the three dimensional structure of a molecule, the improvement comprising means for correlating said three dimensional structure of a G-CSF molecule with the composition of said G-CSF molecule.
- 35 6. A method for crystallization of a protein comprising the steps of:
- (a) combining, optionally by automated means, aqueous aliquots of said protein with either (i) aliquots of a salt solution, each aliquot having a different concentration of salt; or (ii) aliquots of a precipitant solution, each aliquot having a different concentration of precipitant;
 - (b) selecting at least one of said combined aliquots, said selection based on the formation of precrystalline forms, or, if no precrystalline forms are so produced, increasing the protein starting concentration of said aqueous aliquots of protein and repeating step (a);
 - (c) after said salt or said precipitant concentration is selected, repeating step (a) with said previously unselected solution in the presence of said selected concentration; and,
 - (d) repeating step (b) and step (a) until a crystal of desired quality is obtained.
- 40 45 7. A method of claim 6 wherein each combination pursuant to step (a) is performed in a range of pH.
- 50 8. A method of claim 6 wherein said combining of step (a) is done in the presence of a nucleation initiation unit.
9. A G-CSF analog having an amino acid sequence different from that of Figure 1 in that:
- (a) the N-terminal methionine is optional; and
 - (b) one or more of amino acids 58-72 (i) is substituted with one or more different amino acids or (ii) deleted; or (iii) chemically modified.
- 55

10. A G-CSF analog of claim 9 wherein said analog is more resistant to proteolysis than a G-CSF molecule of Figure 1.
11. A G-CSF analog of claim 10 wherein at least one of said amino acids is chemically modified by the addition of a polyethylene glycol molecule.
12. A G-CSF analog having an amino acid sequence different from that of Figure 1 in that:
 - (a) the N-terminal methionine is optional; and
 - (b) one or more of amino acids 119-125 (i) is substituted with one or more different amino acids or (ii) deleted; or (iii) chemically modified.
13. A G-CSF analog of claim 12 wherein said analog is more resistant to proteolysis than a G-CSF molecule of Figure 1.
14. A G-CSF analog of claim 12 wherein at least one of said amino acids is chemically modified by the addition of a polyethylene glycol molecule.
15. A G-CSF molecule having the AB loop stabilized by connecting such loop to one or more of helices A, B, C, or D.
16. A G-CSF molecule having the CD loop stabilized by connecting such loop to one or more of helices A, B, C, or D.
17. A G-CSF analog, optionally in a pharmaceutically effective carrier, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys¹⁷->Arg¹⁷ and the N-terminal methionine is optional.
18. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys³⁵->Arg³⁵ and the N-terminal methionine is optional.
19. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys⁴¹->Arg⁴¹ and the N-terminal methionine is optional.
20. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys^{17,24,35,41}->Arg^{17,24,35,41} and the N-terminal methionine is optional.
21. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys^{17,35,41}->Arg^{17,35,41} and the N-terminal methionine is optional.
22. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys^{24,35,41}->Arg^{24,35,41} and the N-terminal methionine is optional.
23. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys^{17,24,35,41}->Arg^{17,24,35,41} and the N-terminal methionine is optional.
24. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys^{17,24,41}->Arg^{17,24,41} and the N-terminal methionine is optional.
25. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Gln⁶⁸->Glu⁶⁸ and the N-terminal methionine is optional.
26. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Cys^{37,43}->Ser^{37,43} and the N-terminal methionine is optional.
27. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Gln²⁶->Ala²⁶ and the N-terminal methionine is optional.

28. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Gln¹⁷⁴->Ala¹⁷⁴ and the N-terminal methionine is optional.
29. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence
5 differs from that of Figure 1 in that Arg¹⁷⁰->Ala¹⁷⁰ and the N-terminal methionine is optional.
30. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Arg¹⁶⁷->Ala¹⁶⁷ and the N-terminal methionine is optional.
- 10 31. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that there is a deletion at position 167 and the N-terminal methionine is optional.
32. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence
15 differs from that of Figure 1 in that Lys⁴¹->Ala⁴¹ and the N-terminal methionine is optional.
33. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that His⁴⁴->Lys⁴⁴ and the N-terminal methionine is optional.
- 20 34. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Glu⁴⁷->Ala⁴⁷ and the N-terminal methionine is optional.
35. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Arg²³->Ala²³ and the N-terminal methionine is optional.
25
36. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys²⁴->Ala²⁴ and the N-terminal methionine is optional.
37. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence
30 differs from that of Figure 1 in that Glu²⁰->Ala²⁰ and the N-terminal methionine is optional.
38. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Asp²⁸->Ala²⁸ and the N-terminal methionine is optional.
- 35 39. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Met¹²⁷->Glu¹²⁷ and the N-terminal methionine is optional.
40. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from tha of Figure 1 in that Met¹³⁸->Glu¹³⁸ and the N-terminal methionine is optional.
40
41. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Met¹²⁷->Leu¹²⁷ and the N-terminal methionine is optional.
42. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence
45 differs from that of Figure 1 in that Met¹³⁸->Leu¹³⁸ and the N-terminal methionine is optional.
43. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Cys¹⁸->Ala¹⁸ and the N-terminal methionine is optional.
- 50 44. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Gln^{12,21}->Glu^{12,21} and the N-terminal methionine is optional.
45. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence
55 differs from that of Figure 1 in that Gln^{12,21,68}->Glu^{12,21,68} and the N-terminal methionine is optional.
46. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Glu²⁰->Ala²⁰; Ser¹³->Gly¹³ and the N-terminal methionine is optional.

47. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Met^{127,138}->Leu^{127,138} and the N-terminal methionine is optional.
48. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Ser¹³->Ala¹³ and the N-terminal methionine is optional.
49. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Lys¹⁷->Ala¹⁷ and the N-terminal methionine is optional.
50. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Gln¹²¹->Ala¹²¹ and the N-terminal methionine is optional.
51. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Gln²¹->Ala²¹ and the N-terminal methionine is optional.
52. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that His⁴⁴->Ala⁴⁴ and the N-terminal methionine is optional.
53. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein said amino acid sequence differs from that of Figure 1 in that His⁵³->Ala⁵³ and the N-terminal methionine is optional.
54. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Asp¹¹⁰->Ala¹¹⁰ and the N-terminal methionine is optional.
55. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Asp¹¹³->Ala¹¹³ and the N-terminal methionine is optional.
56. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Thr¹¹⁷->Ala¹¹⁷ and the N-terminal methionine is optional.
57. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Asp²⁸->Ala²⁸; Asp¹¹⁰ ->Ala¹¹⁰ and the N-terminal methionine is optional.
58. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Glu¹²⁴->Ala¹²⁴ and the N-terminal methionine is optional.
59. A G-CSF analog, optionally in a pharmaceutically effective carrier, wherein the amino acid sequence differs from that of Figure 1 in that Phe¹¹⁴->Val¹¹⁴, Thr¹¹⁷->A¹¹⁷ and the N-terminal methionine is optional.
60. The G-CSF analog DNA-containing plasmids and bacterial host cells transformed therewith available from the American Type Culture Collection under the accession numbers ATCC 69184, 69185, 69186, 69187, 69188, 69189, 69190, 69191, 69192, 69193, 69194, 69195, 69196, 69197, 69198, 69199, 69200, 69201, 69202, 69203, 69204, 69205, 69206, 69207, 69208, 69209, 69210, 69211, 69212, 69213, 69214, 69215, 69216, 69217, 69218, 69219, 69220, 69221, 69222, 69223, 69224, 69225 and 69226.

Met Thr Pro Leu Gly Pro Ala
TCTAGAAATAACCAAGTAGGTAATAATA ATG ACT CCA TTA GGT CCT CCT

Ser Ser Leu Pro Gln Ser Phe Leu Leu Lys Cys Leu Gly Gln
TCT TCT CTG CCG CAA AGC TTT CTG CTG AAA TGT CTG GAA CAG

Val Arg Lys Ile Gln Gly Asp Gly Ala Ala Leu Gln Glu Lys Leu
GTT CGT AAA ATC CAG GGT GAC GGT GCT GCA CTG CAA GAA AAA CTG

Cys Ala Thr Tyr Lys Leu Cys His Pro Glu Glu Leu Val Leu Leu
TGC GCT ACT TAC AAA CTG TGC CAT CCG GAA GAG CTG GTA CTG CTG

Gly His Ser Leu Gly Ile Pro Trp Ala Pro Leu Ser Ser Cys Pro
GGT CAT TCT CTT GGG ATC CCG TGG GCT CCG CTG TCT TCT TGT CCA

Ser Gln Ala Leu Gln Leu Ala Gly Cys Leu Ser Gln Leu His Ser
TCT CAA GCT CTT CAG CTG GCT GGT TGT CTG TCT CAA CTG CAT TCT

Gly Leu Phe Leu Tyr Gln Gly Leu Leu Gln Ala Leu Glu Gly Ile
GGT CTG TTC CTG TAT CAG GGT CTT CTG CAA GCT CTG GAA GGT ATC

Ser Pro Glu Leu Gly Pro Thr Leu Asp Thr Leu Gln Leu Asp Val
TCT CCG GAA CTG GGT CCG ACT CTG GAC ACT CTG CAG CTA GAT GTA

Ala Asp Phe Ala Thr Thr Ile Trp Gln Gln Met Glu Glu Leu Gly
GCT GAC TTT GCT ACT ACT ATT TGG CAA CAG ATG GAA GAG CTC GGT

Met Ala Pro Ala Leu Gln Pro Thr Gln Gly Ala Met Pro Ala Phe
ATG GCA CCA GCT CTG CAA CCG ACT CAA GGT GCT ATG CCG GCA TTC

Ala Ser Ala Phe Gln Arg Arg Ala Gly Gly Val Leu Val Ala Ser
GCT TCT GCA TTC CAG CGT CGT GCA GGA GGT GTA CTG GTT GCT TCT

His Leu Gln Ser Phe Leu Glu Val Ser Tyr Arg Val Leu Arg His
CAT CTG CAA TCT TTC CTG GAA GTA TCT TAC CGT GTT CTG CGT CAT

Leu Ala Gln Pro OC AM
CTG GCT CAG CCG TAA TAG AATTC

FIGURE 1

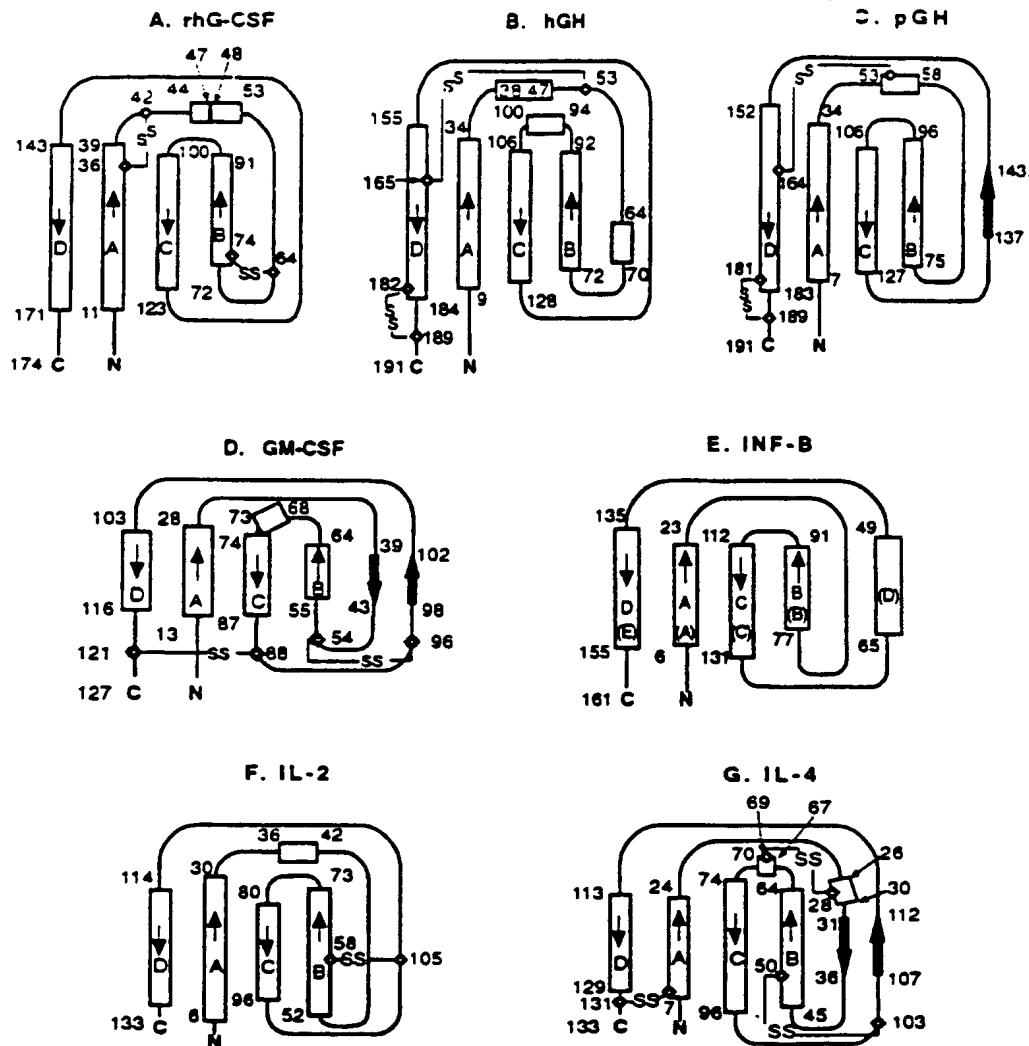


FIGURE 2

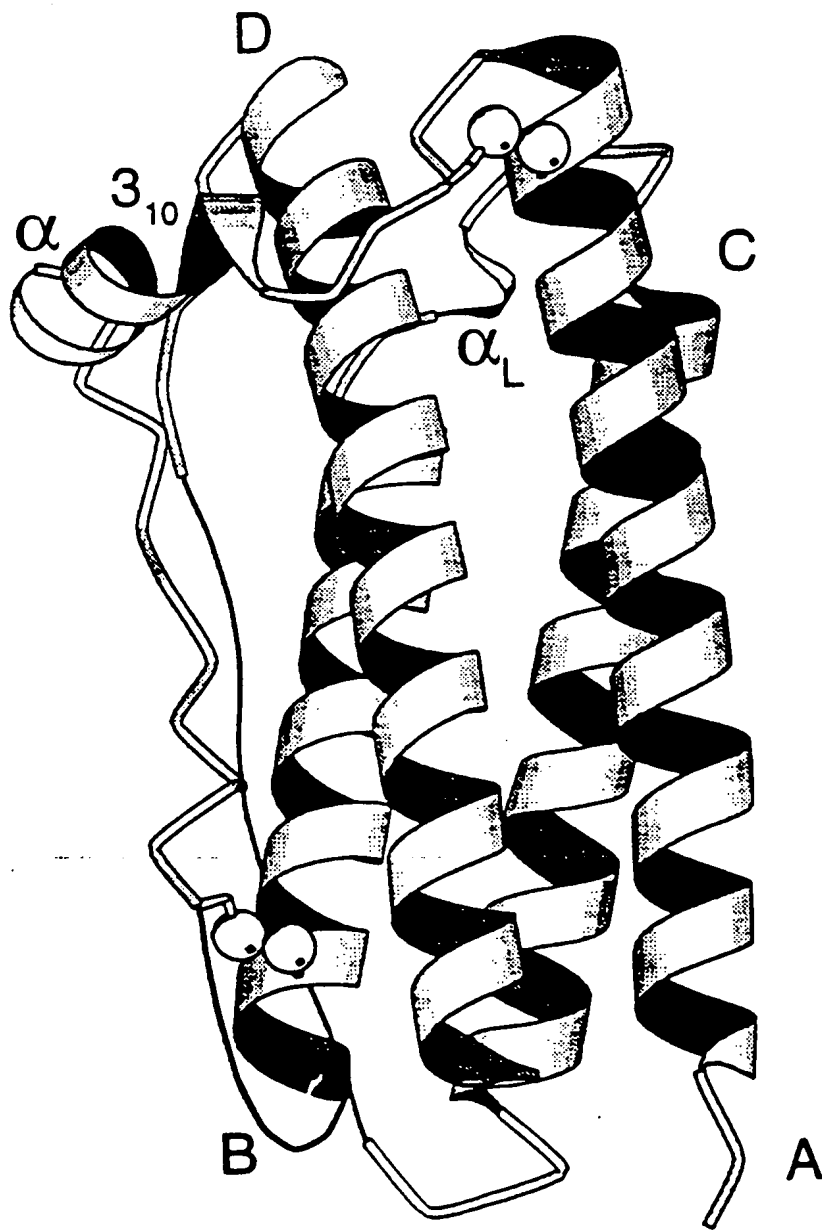


FIGURE 3

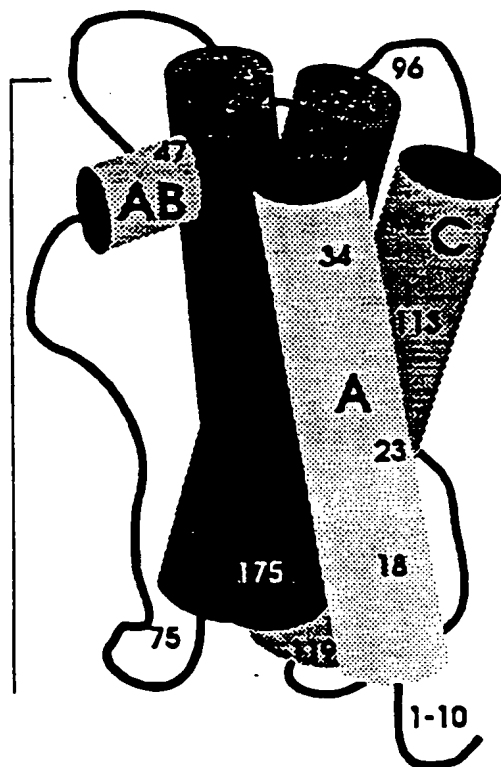


FIGURE 4

FIGURE 5

ATOM	1 CB LEU	10	58.751	58.191	-14.868	1.00	61.22	AI	ATOM	50 O PHE	14	56.789	57.588	-6.805	1.00	46.07	AI
ATOM	2 CG LEU	10	58.360	59.271	-13.939	1.00	60.19	AI	ATOM	51 N LEU	15	57.298	56.509	-8.718	1.00	44.64	AI
ATOM	3 CD1 LEU	10	59.307	60.461	-14.022	1.00	60.14	AI	ATOM	52 N LEU	15	58.024	56.183	-9.287	1.00	41.00	AI
ATOM	4 CD2 LEU	10	56.954	59.658	-14.335	1.00	60.68	AI	ATOM	53 CA LEU	15	55.940	56.181	-9.018	1.00	44.54	AI
ATOM	5 C LEU	10	60.544	56.734	-13.849	1.00	62.85	AI	ATOM	54 CB LEU	15	55.858	55.402	-10.300	1.00	48.72	AI
ATOM	6 O LEU	10	60.079	55.595	-14.041	1.00	63.08	AI	ATOM	55 CG LEU	15	54.853	56.013	-11.269	1.00	51.65	AI
ATOM	7 HT1 LEU	10	59.876	56.135	-15.998	1.00	63.00	AI	ATOM	56 CD1 LEU	15	55.525	57.121	-12.105	1.00	50.34	AI
ATOM	8 HT2 LEU	10	61.323	56.887	-16.434	1.00	62.00	AI	ATOM	57 C112 LEU	15	54.320	54.906	-12.204	1.00	53.77	AI
ATOM	9 N LEU	10	60.328	57.059	-16.704	1.00	62.24	AI	ATOM	58 C LEU	15	55.169	55.410	-8.014	1.00	44.07	AI
ATOM	10 HT3 LEU	10	59.817	57.535	-16.971	1.00	62.58	AI	ATOM	59 O LEU	15	53.945	55.567	-7.959	1.00	45.40	AI
ATOM	11 CA LEU	10	60.183	57.758	-14.941	1.00	62.58	AI	ATOM	60 N LEU	16	55.809	54.620	-7.106	1.00	43.18	AI
ATOM	12 N PRO	11	61.357	56.962	-12.780	1.00	61.96	AI	ATOM	61 H LEU	16	56.781	54.503	-7.251	1.00	61.01	AI
ATOM	13 CD PRO	11	61.960	58.238	-12.383	1.00	61.21	AI	ATOM	62 CA LEU	16	55.110	53.913	-6.095	1.00	42.96	AI
ATOM	14 CA PRO	11	61.832	55.889	-11.906	1.00	61.34	AI	ATOM	63 CB LEU	16	55.866	52.623	-5.751	1.00	11.14	AI
ATOM	15 CB PRO	11	62.915	56.547	-11.043	1.00	59.77	AI	ATOM	64 CG LEU	16	55.840	51.608	-6.868	1.00	42.25	AI
ATOM	16 CG PRO	11	62.511	57.983	-10.975	1.00	59.16	AI	ATOM	65 CD1 LEU	16	56.889	50.567	-6.596	1.00	41.68	AI
ATOM	17 C PRO	11	60.712	55.225	-11.109	1.00	60.68	AI	ATOM	66 CD2 LEU	16	54.413	51.068	-7.030	1.00	42.75	AI
ATOM	18 O PRO	11	60.075	55.843	-10.250	1.00	61.73	AI	ATOM	67 C LEU	16	54.963	54.778	-4.852	1.00	42.35	AI
ATOM	19 N GLN	12	60.466	53.946	-11.407	1.00	59.31	AI	ATOM	68 O LEU	16	54.077	54.579	-4.018	1.00	42.65	AI
ATOM	20 H GLN	12	60.944	53.573	-12.175	1.00	0.00	AI	ATOM	69 N LYS	17	55.821	55.779	-4.703	1.00	42.47	AI
ATOM	21 CA GLN	12	59.468	53.121	-10.743	1.00	57.22	AI	ATOM	70 H LYS	17	56.587	55.840	-5.320	1.00	01.00	AI
ATOM	22 CB GLN	12	59.779	51.646	-10.970	1.00	59.27	AI	ATOM	71 CA LYS	17	55.681	56.767	-3.650	1.00	42.07	AI
ATOM	23 CG GLN	12	58.620	50.714	-10.591	1.00	59.70	AI	ATOM	72 CB LYS	17	56.905	57.554	-3.573	1.00	44.14	AI
ATOM	24 CD GLN	12	57.604	50.575	-11.702	1.00	61.71	AI	ATOM	73 CG LYS	17	57.214	58.197	-2.223	1.00	49.61	AI
ATOM	25 OE1 GLN	12	57.170	49.465	-11.970	1.00	65.82	AI	ATOM	74 CD LYS	17	57.114	57.164	-1.086	1.00	55.15	AI
ATOM	26 NE2 GLN	12	57.227	51.534	-12.541	1.00	63.02	AI	ATOM	75 CE LYS	17	56.747	57.804	0.293	1.00	62.05	AI
ATOM	27 HE21 GLN	12	57.659	52.419	-12.489	1.00	0.00	AI	ATOM	76 NZ1 LYS	17	54.462	58.533	0.331	1.00	65.43	AI
ATOM	28 HE22 GLN	12	56.500	51.308	-13.156	1.00	0.00	AI	ATOM	77 H21 LYS	17	54.684	57.884	0.098	1.00	01.00	AI
ATOM	29 C GLN	12	59.336	53.347	-9.245	1.00	55.34	AI	ATOM	78 H22 LYS	17	55.482	59.308	-0.362	1.00	01.00	AI
ATOM	30 O GLN	12	58.242	53.196	-8.708	1.00	54.56	AI	ATOM	79 H23 LYS	17	55.312	58.926	1.282	1.00	01.00	AI
ATOM	31 N SER	13	60.423	53.732	-8.576	1.00	53.44	AI	ATOM	80 C LYS	17	54.463	57.640	-4.051	1.00	41.70	AI
ATOM	32 H SER	13	61.276	53.839	-9.033	1.00	0.00	AI	ATOM	81 O LYS	17	53.648	57.999	-3.186	1.00	40.66	AI
ATOM	33 CA SER	13	60.335	53.974	-7.168	1.00	52.86	AI	ATOM	82 N CYS	18	54.272	57.992	-5.346	1.00	39.13	AI
ATOM	34 CB SER	13	61.704	54.144	-6.626	1.00	52.24	AI	ATOM	83 H CYS	18	54.798	57.809	-5.981	1.00	01.00	AI
ATOM	35 OG SER	13	61.702	53.493	-5.362	1.00	56.64	AI	ATOM	84 CA CYS	18	53.080	58.656	-5.802	1.00	37.42	AI
ATOM	36 HG SER	13	61.534	52.551	-5.477	1.00	0.00	AI	ATOM	85 CB CYS	18	54.021	60.026	-7.261	1.00	35.02	AI
ATOM	37 C SER	13	59.497	55.214	-6.900	1.00	52.58	AI	ATOM	86 SG CYS	18	54.421	60.026	-7.681	1.00	40.40	AI
ATOM	38 O SER	13	58.509	55.144	-6.160	1.00	53.55	AI	ATOM	87 C CYS	18	51.859	57.789	-5.502	1.00	39.31	AI
ATOM	39 N PHE	14	59.791	56.333	-7.577	1.00	50.84	AI	ATOM	88 O CYS	18	50.959	58.346	-4.847	1.00	40.83	AI
ATOM	40 H PHE	14	60.469	56.292	-8.279	1.00	0.00	AI	ATOM	89 N LEU	19	51.738	56.475	-5.842	1.00	37.15	AI
ATOM	41 CA PHE	14	59.067	57.590	-7.423	1.00	47.21	AI	ATOM	90 H LEU	19	52.462	56.038	-6.341	1.00	01.00	AI
ATOM	42 CB PHE	14	59.611	58.590	-8.454	1.00	44.68	AI	ATOM	91 CA LEU	19	50.571	55.702	-5.534	1.00	36.00	AI
ATOM	43 CG PHE	14	58.618	59.669	-8.866	1.00	42.88	AI	ATOM	92 CB LEU	19	50.644	54.204	-5.917	1.00	38.11	AI
ATOM	44 CD1 PHE	14	58.052	59.594	-10.123	1.00	40.40	AI	ATOM	93 CG LEU	19	49.410	53.271	-5.657	1.00	40.86	AI
ATOM	45 CD2 PHE	14	58.264	60.673	-7.978	1.00	40.40	AI	ATOM	94 CD1 LEU	19	48.208	53.684	-6.467	1.00	39.71	AI
ATOM	46 CE1 PHE	14	57.114	60.518	-10.507	1.00	39.59	AI	ATOM	95 C12 LEU	19	49.602	51.833	-6.113	1.00	45.71	AI
ATOM	47 CE2 PHE	14	57.329	61.587	-8.380	1.00	41.82	AI	ATOM	96 C LEU	19	50.102	55.736	-4.076	1.00	33.52	AI
ATOM	48 CZ PHE	14	56.751	61.515	-9.635	1.00	41.56	AI	ATOM	97 O LEU	19	48.930	55.949	-3.716	1.00	32.75	AI
ATOM	49 C PHE	14	57.605	57.263	-7.661	1.00	45.83	AI	ATOM	98 N GLU	20	51.030	55.576	-3.166	1.00	11.88	AI
ATOM									ATOM	99 H GLU	20	51.940	55.338	-3.455	1.00	61.00	AI
ATOM									ATOM	100 CA GLU	20	50.750	55.710	-1.748	1.00	31.40	AI

FIGURE 5

ATOM	101	CB	GLU	20	52.053	55.334	-1.167	1.00	35.25	AI	ATOM	152	NZ	LYS	24	51.532	59.975	3.333	1.00	51.19	AI
ATOM	102	CG	GLU	20	52.508	55.504	0.760	1.00	43.21	AI	ATOM	153	IZ1	LYS	24	51.637	60.498	4.225	1.00	0.00	AI
ATOM	103	CD	GLU	20	53.948	54.947	0.407	1.00	51.06	AI	ATOM	154	IZ2	LYS	24	51.539	60.651	2.539	1.00	0.00	AI
ATOM	104	OE1	GLU	20	54.370	54.660	1.546	1.00	56.78	AI	ATOM	155	IZ3	LYS	24	52.317	59.303	3.216	1.00	0.00	AI
ATOM	105	OE2	GLU	20	54.708	54.766	-0.570	1.00	51.57	AI	ATOM	156	C	LYS	24	45.455	59.893	1.101	1.00	21.06	AI
ATOM	106	C	GLU	20	50.230	57.117	-1.326	1.00	33.25	AI	ATOM	157	O	LYS	24	44.588	60.068	1.962	1.00	20.90	AI
ATOM	107	O	GLU	20	49.432	57.291	-0.380	1.00	33.30	AI	ATOM	158	N	ILE	25	45.549	60.696	0.044	1.00	21.04	AI
ATOM	108	N	GLN	21	50.660	58.167	-2.044	1.00	32.33	AI	ATOM	159	H	ILE	25	46.242	60.509	-0.629	1.00	0.00	AI
ATOM	109	H	GLN	21	51.270	58.004	-2.794	1.00	0.00	AI	ATOM	160	CA	ILE	25	44.067	61.841	-0.115	1.00	22.53	AI
ATOM	110	CA	GLN	21	50.275	59.538	-1.742	1.00	31.00	AI	ATOM	161	CB	ILE	25	45.075	62.694	-1.307	1.00	22.15	AI
ATOM	111	CB	GLN	21	51.326	60.489	-2.340	1.00	32.37	AI	ATOM	162	CG2	ILE	25	44.097	63.834	-1.439	1.00	20.44	AI
ATOM	112	CG	GLN	21	52.436	60.530	-1.272	1.00	38.01	AI	ATOM	163	CG1	ILE	25	46.475	63.230	-1.136	1.00	21.03	AI
ATOM	113	CD	GLN	21	53.622	61.460	-1.504	1.00	42.67	AI	ATOM	164	CD	ILE	25	47.188	63.281	-2.497	1.00	20.03	AI
ATOM	114	OE1	GLN	21	54.008	62.236	-0.615	1.00	43.63	AI	ATOM	165	C	ILE	25	43.263	61.308	-0.352	1.00	21.75	AI
ATOM	115	NE2	GLN	21	54.256	61.448	-2.678	1.00	42.31	AI	ATOM	166	O	ILE	25	42.339	61.839	0.301	1.00	26.13	AI
ATOM	116	NE21	GLN	21	53.965	60.840	-3.384	1.00	0.00	AI	ATOM	167	N	GLN	26	43.065	60.289	-1.244	1.00	22.79	AI
ATOM	117	NE22	GLN	21	55.026	62.052	-2.730	1.00	0.00	AI	ATOM	168	H	GLN	26	43.842	59.926	-1.726	1.00	0.00	AI
ATOM	118	C	GLN	21	48.894	59.765	-2.288	1.00	28.51	AI	ATOM	169	CA	GLN	26	41.737	59.713	-1.437	1.00	20.12	AI
ATOM	119	O	GLN	21	48.027	60.242	-1.563	1.00	28.65	AI	ATOM	170	CB	GLN	26	41.729	58.539	-2.341	1.00	18.99	AI
ATOM	120	N	VAL	22	48.682	59.319	-3.521	1.00	25.85	AI	ATOM	171	CG	GLN	26	42.203	59.042	-3.627	1.00	19.77	AI
ATOM	121	H	VAL	22	49.448	58.980	-4.013	1.00	0.00	AI	ATOM	172	CD	GLN	26	42.163	57.996	-4.684	1.00	24.26	AI
ATOM	122	CA	VAL	22	47.382	59.303	-4.161	1.00	24.94	AI	ATOM	173	OE1	GLN	26	42.550	56.853	-4.465	1.00	26.82	AI
ATOM	123	CB	VAL	22	47.508	58.614	-5.526	1.00	24.09	AI	ATOM	174	NE2	GLN	26	41.732	58.351	-5.890	1.00	27.08	AI
ATOM	124	CG1	VAL	22	46.154	58.378	-6.096	1.00	19.97	AI	ATOM	175	NE21	GLN	26	41.421	59.265	-6.042	1.00	0.00	AI
ATOM	125	CG2	VAL	22	48.252	59.479	-6.498	1.00	25.82	AI	ATOM	176	NE22	GLN	26	41.743	57.649	-6.552	1.00	0.00	AI
ATOM	126	C	VAL	22	46.418	58.549	-3.226	1.00	25.65	AI	ATOM	177	C	GLN	26	41.707	59.239	-0.111	1.00	21.88	AI
ATOM	127	O	VAL	22	45.428	59.190	-2.800	1.00	29.31	AI	ATOM	178	O	GLN	26	40.067	59.550	0.220	1.00	27.02	AI
ATOM	128	N	ARG	23	46.643	57.291	-2.759	1.00	23.93	AI	ATOM	179	N	GLY	27	41.952	58.622	0.773	1.00	22.54	AI
ATOM	129	H	ARG	23	47.440	56.819	-3.056	1.00	0.00	AI	ATOM	180	H	GLY	27	42.891	58.420	0.575	1.00	0.00	AI
ATOM	130	CA	ARG	23	45.667	56.593	-1.892	1.00	20.67	AI	ATOM	181	CA	GLY	27	41.386	58.191	2.037	1.00	25.55	AI
ATOM	131	CB	ARG	23	46.104	55.135	-1.635	1.00	20.45	AI	ATOM	182	C	GLY	27	40.936	59.352	2.890	1.00	27.80	AI
ATOM	132	CG	ARG	23	46.325	54.321	-2.904	1.00	17.51	AI	ATOM	183	O	GLY	27	39.889	59.251	3.526	1.00	29.95	AI
ATOM	133	CD	ARG	23	45.095	54.446	-3.769	1.00	21.54	AI	ATOM	184	N	ASP	28	41.683	60.460	2.915	1.00	29.49	AI
ATOM	134	NE	ARG	23	45.076	53.437	-4.809	1.00	24.82	AI	ATOM	185	H	ASP	28	42.547	60.454	2.448	1.00	0.00	AI
ATOM	135	HE	ARG	23	45.642	52.647	-4.701	1.00	0.00	AI	ATOM	186	CA	ASP	28	41.257	61.680	3.624	1.00	28.35	AI
ATOM	136	CZ	ARG	23	44.323	53.556	-5.904	1.00	27.69	AI	ATOM	187	CB	ASP	28	42.266	62.789	3.552	1.00	30.11	AI
ATOM	137	NH1	ARG	23	43.562	55.377	-5.303	1.00	0.00	AI	ATOM	188	CG	ASP	28	43.737	62.502	3.777	1.00	31.72	AI
ATOM	138	NH11	ARG	23	43.562	55.377	-5.303	1.00	0.00	AI	ATOM	189	OD1	ASP	28	44.539	63.074	2.995	1.00	31.95	AI
ATOM	139	NH12	ARG	23	42.956	54.730	-6.789	1.00	0.00	AI	ATOM	190	OD2	ASP	28	44.063	61.811	4.241	1.00	32.04	AI
ATOM	140	NH2	ARG	23	44.345	52.604	-6.491	1.00	24.22	AI	ATOM	191	C	ASP	28	39.994	62.264	2.960	1.00	25.81	AI
ATOM	141	HH21	ARG	23	43.780	52.713	-7.709	1.00	0.00	AI	ATOM	192	O	ASP	28	39.101	62.699	3.655	1.00	26.21	AI
ATOM	142	HH22	ARG	23	44.936	51.802	-6.793	1.00	0.00	AI	ATOM	193	N	GLY	29	39.882	62.270	1.631	1.00	23.93	AI
ATOM	143	C	ARG	23	45.458	57.285	-0.560	1.00	20.56	AI	ATOM	194	H	GLY	29	40.660	61.950	1.135	1.00	9.00	AI
ATOM	144	O	ARG	23	44.374	57.254	0.042	1.00	20.04	AI	ATOM	195	CA	GLY	29	38.729	62.694	0.886	1.00	25.69	AI
ATOM	145	N	LYS	24	46.485	58.015	-0.118	1.00	22.67	AI	ATOM	196	C	GLY	29	37.528	61.961	1.418	1.00	27.46	AI
ATOM	146	H	LYS	24	47.291	58.105	-0.668	1.00	0.00	AI	ATOM	197	O	GLY	29	36.648	62.558	2.061	1.00	28.14	AI
ATOM	147	CA	LYS	24	46.431	58.729	1.166	1.00	22.85	AI	ATOM	198	N	ALA	30	37.646	60.628	1.295	1.00	27.85	AI
ATOM	148	CB	LYS	24	47.811	59.255	1.506	1.00	26.86	AI	ATOM	199	H	ALA	30	37.442	60.288	0.843	1.00	0.00	AI
ATOM	149	CD	LYS	24	47.821	59.661	2.971	1.00	33.79	AI	ATOM	200	CA	ALA	30	36.683	59.655	1.814	1.00	25.94	AI
ATOM	150	CE	LYS	24	49.121	60.265	3.404	1.00	40.73	AI	ATOM	201	CB	ALA	30	37.269	58.303	1.556	1.00	22.15	AI
ATOM	151	CE	LYS	24	50.258	59.258	3.335	1.00	46.19	AI	ATOM	202	C	ALA	30	36.356	59.842	3.308	1.00	27.18	AI

FIGURE 5

ATOM 203	O	ALA	30	35.194	59.772	3.754	1.00	28.82	AI	ATOM 254	N	LEU	36	30.452	64.190	6.480	1.00	41.21	AI
ATOM 204	N	ALA	31	37.340	60.105	4.150	1.00	27.16	AI	ATOM 255	H	LEU	36	31.343	61.930	5.836	1.00	0.00	AI
ATOM 205	H	ALA	31	38.253	60.114	3.809	1.00	0.00	AI	ATOM 256	CA	LEU	36	29.647	65.157	6.144	1.00	40.25	AI
ATOM 206	CA	ALA	31	37.113	60.470	5.531	1.00	27.70	AI	ATOM 257	CB	LEU	36	30.070	65.899	4.889	1.00	49.03	AI
ATOM 207	CB	ALA	31	38.383	60.881	6.177	1.00	27.65	AI	ATOM 258	CG	LEU	36	31.253	66.834	4.935	1.00	14.99	AI
ATOM 208	C	ALA	31	36.178	61.675	5.660	1.00	30.01	AI	ATOM 259	CD1	LEU	36	31.438	67.404	3.571	1.00	42.08	AI
ATOM 209	O	ALA	31	35.195	61.624	6.413	1.00	32.91	AI	ATOM 260	CD2	LEU	36	31.034	67.939	5.928	1.00	45.05	AI
ATOM 210	N	LEU	32	36.397	62.744	4.895	1.00	27.63	AI	ATOM 261	C	LEU	36	28.332	64.414	5.941	1.00	41.90	AI
ATOM 211	H	LEU	32	37.133	62.734	4.242	1.00	0.00	AI	ATOM 262	O	LEU	36	27.267	64.828	6.431	1.00	42.10	AI
ATOM 212	CA	LEU	32	35.560	63.898	4.997	1.00	28.52	AI	ATOM 263	N	CYS	37	28.392	63.251	5.309	1.00	42.44	AI
ATOM 213	CB	LEU	32	36.226	65.019	4.167	1.00	32.94	AI	ATOM 264	H	CYS	37	29.250	62.904	5.020	1.00	0.00	AI
ATOM 214	CG	LEU	32	35.658	66.472	4.091	1.00	31.54	AI	ATOM 265	CA	CYS	37	27.216	62.469	5.084	1.00	43.53	AI
ATOM 215	CD1	LEU	32	35.516	67.082	5.499	1.00	32.87	AI	ATOM 266	C	CYS	37	26.638	62.026	6.362	1.00	44.65	AI
ATOM 216	CD2	LEU	32	36.555	67.267	3.181	1.00	30.97	AI	ATOM 267	O	CYS	37	25.426	61.997	6.459	1.00	46.40	AI
ATOM 217	C	LEU	32	34.133	63.597	4.518	1.00	27.87	AI	ATOM 268	CB	CYS	37	27.474	61.240	4.313	1.00	44.00	AI
ATOM 218	O	LEU	32	33.169	63.889	5.350	1.00	25.93	AI	ATOM 269	SG	CYS	37	26.133	60.038	4.530	1.00	43.86	AI
ATOM 219	N	GLN	33	33.977	63.028	3.315	1.00	27.51	AI	ATOM 270	N	ALA	38	27.465	61.734	7.342	1.00	45.96	AI
ATOM 220	H	GLN	33	34.787	62.826	2.802	1.00	0.00	AI	ATOM 271	H	ALA	38	28.433	61.707	7.202	1.00	0.00	AI
ATOM 221	CA	GLN	33	32.687	62.671	2.775	1.00	30.40	AI	ATOM 272	CA	ALA	38	26.932	61.261	8.592	1.00	48.04	AI
ATOM 222	CB	GLN	33	32.737	61.721	1.614	1.00	29.47	AI	ATOM 273	CB	ALA	38	27.869	60.140	9.108	1.00	48.64	AI
ATOM 223	CG	GLN	33	32.888	62.584	0.436	1.00	29.26	AI	ATOM 274	C	ALA	38	26.748	62.358	9.624	1.00	48.91	AI
ATOM 224	CD	GLN	33	33.015	61.869	-0.887	1.00	30.21	AI	ATOM 275	O	ALA	38	26.103	62.085	10.621	1.00	50.72	AI
ATOM 225	OE1	GLN	33	34.064	61.495	-1.452	1.00	29.61	AI	ATOM 276	N	THR	39	27.256	63.590	9.512	1.00	50.66	AI
ATOM 226	NE2	GLN	33	31.823	61.759	-1.426	1.00	31.19	AI	ATOM 277	H	THR	39	27.858	63.780	8.770	1.00	0.00	AI
ATOM 227	HE1	GLN	33	31.781	61.328	-2.302	1.00	0.00	AI	ATOM 278	CA	THR	39	26.976	64.638	10.503	1.00	51.54	AI
ATOM 228	HE2	GLN	33	31.042	62.060	-0.914	1.00	0.00	AI	ATOM 279	CB	THR	39	28.179	65.593	10.690	1.00	51.76	AI
ATOM 229	C	GLN	33	31.849	61.963	3.788	1.00	35.60	AI	ATOM 280	OG1	THR	39	29.294	64.826	11.126	1.00	52.65	AI
ATOM 230	O	GLN	33	30.715	62.416	4.073	1.00	36.49	AI	ATOM 281	HG1	THR	39	29.749	64.481	10.355	1.00	0.00	AI
ATOM 231	N	GLU	34	32.386	60.925	4.438	1.00	39.81	AI	ATOM 282	CG2	THR	39	27.900	66.655	11.729	1.00	51.62	AI
ATOM 232	H	GLU	34	33.340	60.707	4.328	1.00	0.00	AI	ATOM 283	C	THR	39	25.775	65.466	10.037	1.00	52.17	AI
ATOM 233	CA	GLU	34	31.541	60.131	5.304	1.00	43.24	AI	ATOM 284	O	THR	39	24.886	65.882	10.781	1.00	52.15	AI
ATOM 234	CB	GLU	34	32.228	58.792	5.571	1.00	46.46	AI	ATOM 285	N	TYR	40	25.751	65.720	8.738	1.00	52.83	AI
ATOM 235	CG	GLU	34	33.174	58.721	6.624	1.00	55.01	AI	ATOM 286	H	TYR	40	26.420	65.331	8.159	1.00	0.00	AI
ATOM 236	CD	GLU	34	32.777	58.092	7.930	1.00	60.79	AI	ATOM 287	CA	TYR	40	24.729	66.561	8.165	1.00	52.53	AI
ATOM 237	OE1	GLU	34	33.483	57.186	8.412	1.00	63.26	AI	ATOM 288	CB	TYR	40	25.314	67.872	7.696	1.00	52.15	AI
ATOM 238	OE2	GLU	34	31.724	58.504	8.459	1.00	60.44	AI	ATOM 289	CG	TYR	40	26.399	68.458	8.552	1.00	54.11	AI
ATOM 239	C	GLU	34	31.218	60.877	6.564	1.00	43.59	AI	ATOM 290	CD1	TYR	40	27.678	68.341	8.062	1.00	56.50	AI
ATOM 240	O	GLU	34	30.175	60.631	7.161	1.00	44.87	AI	ATOM 291	CE1	TYR	40	28.719	68.934	8.724	1.00	58.28	AI
ATOM 241	N	LYS	35	32.045	61.811	6.998	1.00	44.80	AI	ATOM 292	CD2	TYR	40	26.122	69.144	9.714	1.00	54.86	AI
ATOM 242	H	LYS	35	32.923	61.931	6.569	1.00	0.00	AI	ATOM 293	CE2	TYR	40	27.170	69.746	10.378	1.00	56.20	AI
ATOM 243	CA	LYS	35	31.674	62.634	8.134	1.00	45.43	AI	ATOM 294	C2	TYR	40	28.453	69.642	9.872	1.00	58.26	AI
ATOM 244	CB	LYS	35	32.881	63.364	8.686	1.00	47.67	AI	ATOM 295	OH	TYR	40	29.513	70.310	10.463	1.00	61.00	AI
ATOM 245	CG	LYS	35	33.701	62.414	9.510	1.00	52.75	AI	ATOM 296	HH	TYR	40	30.179	70.443	9.782	1.00	0.00	AI
ATOM 246	CD	LYS	35	35.084	63.021	9.548	1.00	57.55	AI	ATOM 297	C	TYR	40	24.035	65.911	6.981	1.00	51.75	AI
ATOM 247	CE	LYS	35	36.067	62.099	10.238	1.00	60.35	AI	ATOM 298	O	TYR	40	23.662	66.578	6.024	1.00	52.52	AI
ATOM 248	H2	LYS	35	35.810	62.064	11.669	1.00	62.91	AI	ATOM 299	N	LYS	41	23.941	64.600	6.965	1.00	50.54	AI
ATOM 249	H21	LYS	35	34.838	61.733	11.840	1.00	0.00	AI	ATOM 300	H	LYS	41	24.474	64.064	7.583	1.00	0.00	AI
ATOM 250	H22	LYS	35	35.930	63.011	12.078	1.00	0.00	AI	ATOM 301	CA	LYS	41	23.112	63.885	6.029	1.00	50.48	AI
ATOM 251	H23	LYS	35	36.477	61.405	12.119	1.00	0.00	AI	ATOM 302	CB	LYS	41	21.641	63.989	6.540	1.00	50.62	AI
ATOM 252	C	LYS	35	30.630	63.660	7.697	1.00	44.45	AI	ATOM 303	CG	LYS	41	21.387	63.326	7.911	1.00	52.11	AI
ATOM 253	O	LYS	35	29.730	63.999	8.478	1.00	44.61	AI	ATOM 304	CD	LYS	41	20.112	63.878	8.574	1.00	55.54	AI

FIGURE 5

ATOM	305	CE	LVS	41	19.578	63.087	9.820	1.00	58.79	AI	ATOM	356	C	GLU	46	23.181	65.584	-6.937	1.00	-42.36	AI
ATOM	306	NZ	LVS	41	18.374	63.648	10.457	1.00	58.31	AI	ATOM	357	O	GLU	46	22.532	66.223	-7.748	1.00	-41.71	AI
ATOM	307	H21	LVS	41	17.605	63.688	9.757	1.00	0.00	AI	ATOM	358	N	GLU	47	22.919	65.563	-5.654	1.00	-41.96	AI
ATOM	308	H22	LVS	41	18.578	64.607	10.803	1.00	0.00	AI	ATOM	359	H	GLU	47	23.507	65.098	-5.028	1.00	0.00	AI
ATOM	309	H23	LVS	41	18.084	63.043	11.252	1.00	0.00	AI	ATOM	360	CA	GLU	47	21.818	66.301	-5.144	1.00	-41.21	AI
ATOM	310	C	LVS	41	23.251	64.318	4.588	1.00	49.92	AI	ATOM	361	CB	GLU	47	21.294	65.487	-3.963	1.00	-41.24	AI
ATOM	311	O	LVS	41	22.312	64.124	3.793	1.00	51.49	AI	ATOM	362	CG	GLU	47	21.409	65.925	-2.515	1.00	-46.07	AI
ATOM	312	N	LEU	42	24.432	64.893	4.246	1.00	48.78	AI	ATOM	363	CD	GLU	47	20.812	64.907	-1.547	1.00	-47.86	AI
ATOM	313	H	LEU	42	25.103	65.050	4.937	1.00	0.00	AI	ATOM	364	OE1	GLU	47	19.847	64.225	-1.910	1.00	50.99	AI
ATOM	314	CA	LEU	42	24.742	65.286	2.859	1.00	46.61	AI	ATOM	365	OE2	GLU	47	21.313	64.780	-0.427	1.00	-49.47	AI
ATOM	315	CB	LEU	42	25.565	66.574	2.757	1.00	44.69	AI	ATOM	366	C	GLU	47	22.295	67.218	-4.809	1.00	-44.04	AI
ATOM	316	CG	LEU	42	24.807	67.802	3.218	1.00	42.63	AI	ATOM	367	O	GLU	47	21.532	68.547	-4.292	1.00	-44.60	AI
ATOM	317	CD1	LEU	42	25.718	68.580	4.097	1.00	43.29	AI	ATOM	368	N	LEU	48	23.567	68.015	-5.121	1.00	-43.05	AI
ATOM	318	CD2	LEU	42	24.283	68.590	2.045	1.00	41.26	AI	ATOM	369	H	LEU	48	24.140	67.310	-5.465	1.00	0.00	AI
ATOM	319	C	LEU	42	25.580	64.124	2.397	1.00	45.46	AI	ATOM	370	CA	LEU	48	24.166	69.318	-4.904	1.00	-42.42	AI
ATOM	320	O	LEU	42	26.768	64.017	2.711	1.00	46.32	AI	ATOM	371	CB	LEU	48	25.223	69.201	-3.858	1.00	-40.53	AI
ATOM	321	N	CYS	43	24.882	63.193	1.754	1.00	44.09	AI	ATOM	372	CG	LEU	48	24.920	68.695	-2.489	1.00	-41.87	AI
ATOM	322	H	CYS	43	23.925	63.353	1.619	1.00	0.00	AI	ATOM	373	CD1	LEU	48	26.277	68.424	-1.892	1.00	-41.71	AI
ATOM	323	CA	CYS	43	25.448	61.846	-0.123	1.00	41.62	AI	ATOM	374	CD2	LEU	48	24.096	69.670	-1.633	1.00	-41.13	AI
ATOM	324	C	CYS	43	25.448	61.846	-0.123	1.00	41.62	AI	ATOM	375	C	LEU	48	24.792	69.937	-6.166	1.00	-42.37	AI
ATOM	325	O	CYS	43	25.762	60.805	-0.666	1.00	41.99	AI	ATOM	376	O	LEU	48	25.439	70.994	-6.098	1.00	-42.37	AI
ATOM	326	CB	CYS	43	24.716	60.796	2.026	1.00	41.77	AI	ATOM	377	N	VAL	49	24.566	69.366	-7.347	1.00	-41.52	AI
ATOM	327	SG	CYS	43	24.523	61.011	3.835	1.00	45.91	AI	ATOM	378	H	VAL	49	23.951	68.602	-7.362	1.00	0.00	AI
ATOM	328	N	IIS	44	25.057	62.846	-0.867	1.00	42.90	AI	ATOM	379	CA	VAL	49	25.191	69.822	-8.578	1.00	-43.44	AI
ATOM	329	H	IIS	44	24.841	63.721	-0.491	1.00	0.00	AI	ATOM	380	CB	VAL	49	24.890	68.761	-9.636	1.00	-44.29	AI
ATOM	330	CA	IIS	44	25.069	62.680	-2.320	1.00	44.60	AI	ATOM	381	CG1	VAL	49	23.381	68.709	-9.830	1.00	-47.50	AI
ATOM	331	CB	IIS	44	23.653	62.264	-2.825	1.00	48.40	AI	ATOM	382	CG2	VAL	49	25.540	69.086	-10.975	1.00	-45.25	AI
ATOM	332	CG	IIS	44	23.085	60.935	-2.310	1.00	50.37	AI	ATOM	383	C	VAL	49	24.740	71.214	-9.028	1.00	-44.98	AI
ATOM	333	CD1	IIS	44	22.178	60.844	-1.272	1.00	50.52	AI	ATOM	384	O	VAL	49	25.401	71.901	-9.814	1.00	-46.03	AI
ATOM	334	ND1	IIS	44	23.358	59.689	-2.713	1.00	52.28	AI	ATOM	385	N	LEU	50	23.565	71.602	-8.530	1.00	-46.16	AI
ATOM	335	HDI	IIS	44	24.130	59.394	-3.251	1.00	0.00	AI	ATOM	386	H	LEU	50	23.081	70.933	-8.006	1.00	0.00	AI
ATOM	336	CE1	IIS	44	22.652	58.873	-1.955	1.00	51.92	AI	ATOM	387	CA	LEU	50	22.908	72.895	-8.729	1.00	-46.03	AI
ATOM	337	NE2	IIS	44	21.947	59.565	-1.091	1.00	50.53	AI	ATOM	388	CB	LEU	50	21.469	72.769	-8.264	1.00	-46.43	AI
ATOM	338	HE2	IIS	44	21.290	59.189	-0.466	1.00	0.00	AI	ATOM	389	CG	LEU	50	20.443	73.718	-8.760	1.00	-44.16	AI
ATOM	339	C	IIS	44	25.522	63.941	-3.047	1.00	43.69	AI	ATOM	390	CD1	LEU	50	20.259	73.558	-10.243	1.00	-44.79	AI
ATOM	340	O	IIS	44	24.765	64.906	-3.108	1.00	43.00	AI	ATOM	391	CD2	LEU	50	19.159	73.400	-8.079	1.00	-44.04	AI
ATOM	341	N	PRO	45	26.710	63.978	-3.667	1.00	43.07	AI	ATOM	392	C	LEU	50	23.632	73.968	-7.917	1.00	-45.85	AI
ATOM	342	CD	PRO	45	27.785	62.995	-3.501	1.00	42.17	AI	ATOM	393	O	LEU	50	23.996	74.989	-8.484	1.00	-44.52	AI
ATOM	343	CA	PRO	45	27.133	65.074	-4.570	1.00	42.50	AI	ATOM	394	N	LEU	51	23.853	73.764	-6.606	1.00	-45.44	AI
ATOM	344	CB	PRO	45	28.380	64.466	-5.217	1.00	39.76	AI	ATOM	395	H	LEU	51	23.489	72.958	-6.189	1.00	0.00	AI
ATOM	345	CG	PRO	45	28.995	63.680	-4.123	1.00	39.09	AI	ATOM	396	CA	LEU	51	24.676	74.656	-5.805	1.00	-46.04	AI
ATOM	346	C	PRO	45	26.071	65.423	-5.585	1.00	44.49	AI	ATOM	397	CB	LEU	51	24.860	74.084	-4.435	1.00	-45.53	AI
ATOM	347	O	PRO	45	25.876	66.612	-5.801	1.00	45.36	AI	ATOM	398	CG	LEU	51	25.741	74.931	-3.535	1.00	-47.78	AI
ATOM	348	N	GLU	46	25.334	64.501	-6.225	1.00	45.36	AI	ATOM	399	CD1	LEU	51	25.148	76.320	-3.322	1.00	-47.33	AI
ATOM	349	H	GLU	46	25.464	63.561	-5.996	1.00	0.00	AI	ATOM	400	CD2	LEU	51	25.902	74.202	-2.219	1.00	-48.33	AI
ATOM	350	CA	GLU	46	24.406	64.806	-7.319	1.00	45.46	AI	ATOM	401	C	LEU	51	26.064	74.845	-6.416	1.00	-46.27	AI
ATOM	351	CB	GLU	46	23.952	63.515	-7.997	1.00	50.54	AI	ATOM	402	O	LEU	51	26.551	75.966	-6.612	1.00	-47.62	AI
ATOM	352	CG	GLU	46	24.467	63.460	-9.445	1.00	58.48	AI	ATOM	403	N	GLY	52	26.702	73.736	-6.809	1.00	-44.84	AI
ATOM	353	CD	GLU	46	23.637	64.215	-10.502	1.00	64.94	AI	ATOM	404	H	GLY	52	26.306	72.869	-6.578	1.00	0.00	AI
ATOM	354	OE1	GLU	46	23.642	65.455	-10.516	1.00	68.55	AI	ATOM	405	CA	GLY	52	27.989	73.758	-7.453	1.00	-47.91	AI
ATOM	355	OE2	GLU	46	22.995	63.554	-11.332	1.00	68.31	AI	ATOM	406	C	GLY	52	27.984	74.533	-8.750	1.00	-42.47	AI

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ATOM	407 O GLY	52	28.853	75.364	-8.983	1.00	42.06	AI	ATOM	458 O PRO	58	37.187	73.599	-9.691	1.00	49.75	AI
ATOM	408 N HIS	53	27.047	74.307	-9.653	1.00	42.02	AI	ATOM	459 N TRP	59	37.030	72.927	-11.816	1.00	50.37	AI
ATOM	409 H HIS	53	26.366	73.624	-9.471	1.00	0.00	AI	ATOM	460 H TRP	59	36.888	73.141	-12.760	1.00	0.00	AI
ATOM	410 CA HIS	53	27.009	75.104	-10.861	1.00	42.23	AI	ATOM	461 CA TRP	59	37.524	71.595	-11.482	1.00	51.78	AI
ATOM	411 CB HIS	53	25.842	74.689	-11.706	1.00	42.21	AI	ATOM	462 CB TRP	59	36.435	70.562	-11.857	1.00	49.06	AI
ATOM	412 CG HIS	53	26.076	73.399	-12.460	1.00	44.60	AI	ATOM	463 CG TRP	59	35.254	70.712	-10.889	1.00	46.37	AI
ATOM	413 CD2 HIS	53	25.112	72.774	-13.200	1.00	47.49	AI	ATOM	464 CD2 TRP	59	35.320	70.845	-9.521	1.00	44.06	AI
ATOM	414 ND1 HIS	53	27.180	72.669	-12.578	1.00	46.76	AI	ATOM	465 CD2 TRP	59	33.908	71.027	-9.205	1.00	44.18	AI
ATOM	415 HD1 HIS	53	28.039	72.853	-12.139	1.00	0.00	AI	ATOM	466 CD2 TRP	59	36.274	70.842	-8.538	1.00	44.03	AI
ATOM	416 CE1 HIS	53	26.954	71.641	-13.346	1.00	46.90	AI	ATOM	467 CD1 TRP	59	33.729	70.794	-11.354	1.00	45.17	AI
ATOM	417 NE2 HIS	53	25.704	71.725	-13.707	1.00	50.22	AI	ATOM	468 NE1 TRP	59	33.729	70.794	-10.297	1.00	43.17	AI
ATOM	418 HE2 HIS	53	25.237	71.033	-14.239	1.00	0.00	AI	ATOM	469 HE1 TRP	59	32.301	71.312	-10.332	1.00	0.00	AI
ATOM	419 C HIS	53	26.893	76.585	-10.536	1.00	42.72	AI	ATOM	470 C22 TRP	59	33.598	71.215	-7.916	1.00	45.60	AI
ATOM	420 O HIS	53	27.622	77.399	-11.068	1.00	42.03	AI	ATOM	471 C23 TRP	59	35.893	71.028	-7.243	1.00	45.25	AI
ATOM	421 N SER	54	26.099	76.920	-9.535	1.00	45.08	AI	ATOM	472 CH2 TRP	59	34.565	71.214	-6.938	1.00	46.44	AI
ATOM	422 H SER	54	23.673	76.218	-9.001	1.00	0.00	AI	ATOM	473 C TRP	59	38.815	71.435	-12.236	1.00	52.84	AI
ATOM	423 CA SER	54	25.792	78.278	-9.177	1.00	46.92	AI	ATOM	474 O TRP	59	38.842	71.972	-13.372	1.00	54.96	AI
ATOM	424 CB SER	54	24.576	78.181	-8.289	1.00	48.86	AI	ATOM	475 N AIA	60	39.912	70.834	-11.777	1.00	51.97	AI
ATOM	425 CG SER	54	23.521	77.616	-9.112	1.00	53.06	AI	ATOM	476 H AIA	60	39.857	70.269	-10.977	1.00	0.00	AI
ATOM	426 HG SER	54	23.465	76.677	-8.918	1.00	0.00	AI	ATOM	477 CA AIA	60	41.108	70.870	-12.609	1.00	52.18	AI
ATOM	427 C SER	54	26.939	79.033	-8.549	1.00	47.92	AI	ATOM	478 CB AIA	60	42.303	70.610	-11.748	1.00	51.75	AI
ATOM	428 O SER	54	27.038	80.264	-8.655	1.00	49.60	AI	ATOM	479 C AIA	60	41.055	69.857	-13.246	1.00	54.16	AI
ATOM	429 N LEU	55	27.837	78.273	-7.933	1.00	47.59	AI	ATOM	480 O AIA	60	40.545	68.760	-13.530	1.00	51.17	AI
ATOM	430 H LEU	55	27.638	77.322	-7.791	1.00	0.00	AI	ATOM	481 N PRO	61	41.435	70.145	-14.986	1.00	53.44	AI
ATOM	431 CA LEU	55	29.075	78.810	-7.401	1.00	45.27	AI	ATOM	482 CD PRO	61	41.370	71.458	-15.622	1.00	54.76	AI
ATOM	432 CB LEU	55	29.552	77.913	-6.243	1.00	45.49	AI	ATOM	483 CA PRO	61	41.691	69.145	-15.993	1.00	55.57	AI
ATOM	433 CG LEU	55	28.840	77.992	-4.874	1.00	47.30	AI	ATOM	484 CB PRO	61	41.792	69.918	-17.310	1.00	54.95	AI
ATOM	434 CD1 LEU	55	28.876	76.596	-4.299	1.00	49.52	AI	ATOM	485 CG PRO	61	42.211	71.297	-16.901	1.00	54.05	AI
ATOM	435 CD2 LEU	55	29.530	78.921	-3.862	1.00	45.69	AI	ATOM	486 C PRO	61	42.934	68.333	-15.690	1.00	57.54	AI
ATOM	436 C LEU	55	30.133	78.889	-8.492	1.00	43.63	AI	ATOM	487 O PRO	61	43.757	68.661	-14.834	1.00	57.20	AI
ATOM	437 O LEU	55	31.247	79.350	-8.272	1.00	43.24	AI	ATOM	488 N LEU	62	43.040	67.271	-16.486	1.00	59.98	AI
ATOM	438 N GLY	56	29.855	78.383	-9.675	1.00	43.55	AI	ATOM	489 H LEU	62	42.285	67.067	-17.077	1.00	0.00	AI
ATOM	439 H GLY	56	28.984	77.975	-9.828	1.00	0.00	AI	ATOM	490 CA LEU	62	44.184	66.370	-16.471	1.00	63.64	AI
ATOM	440 CA GLY	56	30.814	78.590	-10.753	1.00	45.39	AI	ATOM	491 CB LEU	62	44.062	65.417	-15.260	1.00	63.72	AI
ATOM	441 C GLY	56	32.182	77.811	-10.392	1.00	46.76	AI	ATOM	492 CG LEU	62	45.323	64.691	-14.865	1.00	64.43	AI
ATOM	442 O GLY	56	33.171	78.213	-11.015	1.00	47.31	AI	ATOM	493 CD1 LEU	62	46.394	65.704	-14.488	1.00	64.02	AI
ATOM	443 N ILE	57	32.247	76.885	-9.412	1.00	47.49	AI	ATOM	494 CD2 LEU	62	45.016	63.764	-13.717	1.00	64.98	AI
ATOM	444 H ILE	57	31.392	76.594	-9.042	1.00	0.00	AI	ATOM	495 C LEU	62	44.214	65.611	-17.812	1.00	65.69	AI
ATOM	445 CA ILE	57	33.406	76.249	-8.950	1.00	48.28	AI	ATOM	496 OT1 LEU	62	44.256	66.302	-18.844	1.00	68.47	AI
ATOM	446 CB ILE	57	33.144	75.172	-7.863	1.00	47.79	AI	ATOM	497 OT2 LEU	62	44.194	64.371	-17.845	1.00	66.57	AI
ATOM	447 CG2 ILE	57	34.457	74.591	-7.348	1.00	46.85	AI	ATOM	498 CB LEU	72	57.448	63.159	-19.422	1.00	61.44	A2
ATOM	448 CG1 ILE	57	32.338	75.764	-6.701	1.00	45.09	AI	ATOM	499 CG LEU	72	57.716	62.495	-18.117	1.00	61.40	A2
ATOM	449 CD1 ILE	57	31.859	74.739	-5.659	1.00	41.23	AI	ATOM	500 CD1 LEU	72	56.719	61.408	-17.913	1.00	61.50	A2
ATOM	450 C ILE	57	34.276	75.602	-10.115	1.00	49.15	AI	ATOM	501 CD2 LEU	72	55.107	61.901	-18.121	1.00	63.22	A2
ATOM	451 O ILE	57	33.678	74.935	-10.968	1.00	49.04	AI	ATOM	502 C LEU	72	55.897	65.084	-18.876	1.00	65.40	A2
ATOM	452 N PRO	58	35.596	75.817	-10.248	1.00	49.75	AI	ATOM	503 O LEU	72	55.469	65.301	-18.316	1.00	67.30	A2
ATOM	453 CD PRO	58	36.402	76.743	-9.433	1.00	50.94	AI	ATOM	504 HT1 LEU	72	56.469	64.683	-21.761	1.00	0.00	A2
ATOM	454 CA PRO	58	36.421	75.228	-11.302	1.00	50.72	AI	ATOM	505 HT2 LEU	72	54.827	64.355	-20.951	1.00	9.00	A2
ATOM	455 CB PRO	58	37.525	76.241	-11.488	1.00	50.92	AI	ATOM	506 N LEU	72	55.795	63.983	-20.899	1.00	66.79	A2
ATOM	456 CG PRO	58	37.814	76.663	-10.041	1.00	50.82	AI	ATOM	507 HT3 LEU	72	55.866	63.098	-21.419	1.00	0.00	A2
ATOM	457 C PRO	58	36.916	73.845	-10.875	1.00	50.36	AI	ATOM	508 CA LEU	72	56.064	63.714	-19.512	1.00	64.91	A2

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ATOM	509 N	ALA	73	56.807	66.046	-19.086	1.00	64.54	A2	ATOM	560 CG	LEU	79	49.154	64.895	-11.251	1.00	45.18	A2
ATOM	510 H	ALA	73	57.690	65.804	-19.432	1.00	0.00	A2	ATOM	561 CD1	LEU	79	49.634	63.594	-11.957	1.00	45.06	A2
ATOM	511 C	ALA	73	56.707	67.433	-18.615	1.00	62.55	A2	ATOM	562 CD2	LEU	79	49.766	64.986	-9.969	1.00	46.03	A2
ATOM	512 CB	ALA	73	57.553	68.314	-19.529	1.00	64.84	A2	ATOM	563 C	LEU	79	49.366	68.265	-11.170	1.00	44.49	A2
ATOM	513 C	ALA	73	55.319	68.024	-18.539	1.00	60.37	A2	ATOM	564 O	LEU	79	48.643	68.509	-10.199	1.00	43.20	A2
ATOM	514 O	ALA	73	54.801	68.180	-17.456	1.00	59.42	A2	ATOM	565 N	HIS	80	50.556	68.834	-11.329	1.00	43.83	A2
ATOM	515 N	GLY	74	54.693	68.226	-19.691	1.00	59.72	A2	ATOM	566 H	HIS	80	51.115	68.548	-12.085	1.00	0.00	A2
ATOM	516 H	GLY	74	55.212	68.174	-20.514	1.00	0.00	A2	ATOM	567 CA	HIS	80	51.060	69.788	-10.360	1.00	44.79	A2
ATOM	517 CA	GLY	74	53.336	68.728	-19.816	1.00	59.99	A2	ATOM	568 CG	HIS	80	52.456	70.221	-10.810	1.00	43.58	A2
ATOM	518 C	GLY	74	52.327	68.114	-18.865	1.00	60.27	A2	ATOM	569 CB	HIS	80	53.030	71.031	-9.690	1.00	43.75	A2
ATOM	519 O	GLY	74	51.880	68.796	-17.935	1.00	60.80	A2	ATOM	570 CD2	HIS	80	53.484	70.497	-8.517	1.00	47.48	A2
ATOM	520 N	CYS	75	51.945	66.850	-19.030	1.00	59.60	A2	ATOM	571 ND1	HIS	80	53.083	72.343	-9.567	1.00	44.74	A2
ATOM	521 H	CYS	75	52.160	66.358	-19.839	1.00	0.00	A2	ATOM	572 HD1	HIS	80	52.842	73.004	-10.255	1.00	0.00	A2
ATOM	522 CA	CYS	75	51.002	66.276	-18.078	1.00	60.17	A2	ATOM	573 CE1	HIS	80	53.530	72.641	-8.376	1.00	44.47	A2
ATOM	523 CB	CYS	75	50.670	64.801	-18.464	1.00	64.08	A2	ATOM	574 NE2	HIS	80	53.772	71.520	-7.748	1.00	48.16	A2
ATOM	524 SG	CYS	75	49.832	64.732	-20.096	1.00	73.47	A2	ATOM	575 HE2	HIS	80	54.103	71.444	-6.824	1.00	0.00	A2
ATOM	525 C	CYS	75	51.502	66.346	-16.642	1.00	56.73	A2	ATOM	576 C	HIS	80	50.094	70.978	-10.219	1.00	44.40	A2
ATOM	526 O	CYS	75	50.734	66.748	-15.765	1.00	55.82	A2	ATOM	577 O	HIS	80	49.643	71.294	-9.131	1.00	44.28	A2
ATOM	527 N	LEU	76	52.795	66.142	-16.396	1.00	53.93	A2	ATOM	578 N	SER	81	49.733	71.670	-11.309	1.00	45.13	A2
ATOM	528 H	LEU	76	53.423	66.043	-17.137	1.00	0.00	A2	ATOM	579 H	SER	81	50.136	71.459	-12.176	1.00	0.00	A2
ATOM	529 CA	LEU	76	53.325	66.156	-15.044	1.00	52.94	A2	ATOM	580 CA	SER	81	48.738	72.742	-11.296	1.00	45.41	A2
ATOM	530 CB	LEU	76	54.798	65.754	-15.181	1.00	50.81	A2	ATOM	581 CB	SER	81	48.612	73.347	-12.682	1.00	45.59	A2
ATOM	531 CG	LEU	76	55.575	65.011	-14.090	1.00	49.02	A2	ATOM	582 CG	SER	81	49.594	73.444	-13.292	1.00	49.27	A2
ATOM	532 CD1	LEU	76	54.852	63.740	-13.698	1.00	46.76	A2	ATOM	583 HG	SER	81	50.058	72.670	-13.843	1.00	0.00	A2
ATOM	533 CD2	LEU	76	56.951	64.633	-14.623	1.00	47.67	A2	ATOM	584 C	SER	81	47.344	72.266	-10.855	1.00	44.85	A2
ATOM	534 C	LEU	76	53.093	67.545	-14.425	1.00	53.65	A2	ATOM	585 O	SER	81	46.604	73.064	-10.256	1.00	46.83	A2
ATOM	535 O	LEU	76	52.731	67.716	-13.244	1.00	53.50	A2	ATOM	586 N	GLY	82	46.946	71.010	-11.092	1.00	42.16	A2
ATOM	536 N	SER	77	53.137	68.553	-15.301	1.00	53.91	A2	ATOM	587 H	GLY	82	47.513	70.411	-11.614	1.00	0.00	A2
ATOM	537 H	SER	77	53.322	68.361	-16.242	1.00	0.00	A2	ATOM	588 CA	GLY	82	45.663	70.500	-10.650	1.00	59.59	A2
ATOM	538 CA	SER	77	52.882	69.932	-14.942	1.00	54.93	A2	ATOM	589 C	GLY	82	45.569	70.461	-9.139	1.00	59.50	A2
ATOM	539 CB	SER	77	53.425	70.835	-16.040	1.00	58.32	A2	ATOM	590 O	GLY	82	44.542	70.843	-8.544	1.00	59.64	A2
ATOM	540 CG	SER	77	54.806	70.387	-16.310	1.00	63.35	A2	ATOM	591 N	LEU	83	46.576	70.032	-8.521	1.00	37.57	A2
ATOM	541 HG	SER	77	54.949	69.637	-16.315	1.00	0.00	A2	ATOM	592 H	LEU	83	47.413	69.695	-9.075	1.00	0.00	A2
ATOM	542 C	SER	77	51.382	70.172	-14.759	1.00	53.47	A2	ATOM	593 CA	LEU	83	46.826	70.007	-7.057	1.00	38.07	A2
ATOM	543 O	SER	77	50.982	70.965	-13.899	1.00	53.54	A2	ATOM	594 CB	LEU	83	46.133	69.202	-6.748	1.00	35.67	A2
ATOM	544 N	GLN	78	50.509	69.501	-13.512	1.00	51.82	A2	ATOM	595 CG	LEU	83	48.071	67.736	-7.245	1.00	32.51	A2
ATOM	545 H	GLN	78	50.857	68.901	-16.207	1.00	0.00	A2	ATOM	596 CD1	LEU	83	49.442	67.145	-7.319	1.00	29.77	A2
ATOM	546 CA	GLN	78	49.074	69.619	-15.349	1.00	50.74	A2	ATOM	597 CD2	LEU	83	47.180	66.973	-6.288	1.00	28.71	A2
ATOM	547 CB	GLN	78	48.402	68.877	-16.451	1.00	54.31	A2	ATOM	598 C	LEU	83	46.836	71.386	-6.354	1.00	38.48	A2
ATOM	548 CG	GLN	78	47.420	69.784	-17.160	1.00	58.59	A2	ATOM	599 O	LEU	83	46.392	71.627	-5.219	1.00	38.05	A2
ATOM	549 CD	GLN	78	46.557	68.940	-18.071	1.00	62.32	A2	ATOM	600 N	PIE	84	47.366	72.338	-7.108	1.00	40.34	A2
ATOM	550 OE1	GLN	78	47.005	68.260	-18.998	1.00	65.94	A2	ATOM	601 H	PIE	84	47.804	72.078	-7.944	1.00	0.00	A2
ATOM	551 NE2	GLN	78	45.269	68.889	-17.800	1.00	63.17	A2	ATOM	602 CA	PIE	84	47.414	73.703	-6.088	1.00	41.54	A2
ATOM	552 HE21	GLN	78	44.973	69.327	-16.972	1.00	0.00	A2	ATOM	603 CB	PIE	84	48.163	74.531	-7.693	1.00	46.88	A2
ATOM	553 HE22	GLN	78	44.704	68.444	-18.456	1.00	0.00	A2	ATOM	604 CG	PIE	84	48.715	75.777	-6.988	1.00	55.09	A2
ATOM	554 C	GLN	78	48.591	69.065	-14.011	1.00	48.17	A2	ATOM	605 CD1	PIE	84	49.521	75.622	-5.849	1.00	55.41	A2
ATOM	555 O	GLN	78	47.691	69.618	-13.368	1.00	46.31	A2	ATOM	606 CD2	PIE	84	48.596	77.053	-7.469	1.00	55.79	A2
ATOM	556 N	LEU	79	49.236	67.988	-13.564	1.00	45.89	A2	ATOM	607 CE1	PIE	84	50.004	76.737	-5.195	1.00	57.60	A2
ATOM	557 H	LEU	79	49.920	67.584	-14.140	1.00	0.00	A2	ATOM	608 CE2	PIE	84	48.392	78.156	-6.796	1.00	57.25	A2
ATOM	558 CA	LEU	79	48.919	67.359	-12.294	1.00	44.54	A2	ATOM	609 CZ	PIE	84	49.688	78.092	-5.667	1.00	58.44	A2
ATOM	559 CB	LEU	79	49.617	66.015	-12.259	1.00	45.06	A2	ATOM	610 C	PIE	84	45.994	74.191	-6.531	1.00	40.47	A2

FIGURE 5

ATOM	611	O	PIUE	84	45.609	74.749	-5.558	1.00	42.71	A2	ATOM	662	II	LEU	90	41.101	73.626	-1.643	1.00	0.00	A2
ATOM	612	N	LEU	85	45.190	73.953	-7.624	1.00	38.64	A2	ATOM	663	CA	LEU	90	40.182	73.274	0.235	1.00	51.41	A2
ATOM	613	II	LEU	85	45.555	73.527	-8.429	1.00	0.00	A2	ATOM	664	CB	LEU	90	41.207	72.234	0.501	1.00	36.15	A2
ATOM	614	CA	LEU	85	43.794	74.335	-7.584	1.00	38.81	A2	ATOM	665	CG	LEU	90	41.075	70.971	-0.343	1.00	38.76	A2
ATOM	615	CB	LEU	85	43.101	73.886	-8.839	1.00	41.27	A2	ATOM	666	CD1	LEU	90	42.431	70.267	-0.456	1.00	37.21	A2
ATOM	616	CG	LEU	85	41.673	74.403	-9.017	1.00	46.45	A2	ATOM	667	CD2	LEU	90	39.995	70.099	0.279	1.00	40.54	A2
ATOM	617	CD1	LEU	85	41.702	75.784	-9.719	1.00	47.80	A2	ATOM	668	C	LEU	90	40.342	74.319	1.255	1.00	34.21	A2
ATOM	618	CD2	LEU	85	40.860	73.359	-9.787	1.00	48.25	A2	ATOM	669	O	LEU	90	39.711	74.256	2.313	1.00	35.57	A2
ATOM	619	C	LEU	85	43.079	73.731	-6.386	1.00	38.20	A2	ATOM	670	N	GIN	91	41.188	75.291	0.940	1.00	35.24	A2
ATOM	620	O	LEU	85	42.498	74.469	-5.582	1.00	38.36	A2	ATOM	671	II	GIN	91	41.563	75.284	0.078	1.00	0.00	A2
ATOM	621	N	TYR	86	43.150	72.405	-6.198	1.00	37.92	A2	ATOM	672	CA	GIN	91	41.397	76.373	1.883	1.00	37.40	A2
ATOM	622	N	TYR	86	43.637	71.850	-6.845	1.00	0.00	A2	ATOM	673	CB	GIN	91	42.557	77.182	1.363	1.00	39.65	A2
ATOM	623	CA	TYR	86	42.501	71.801	-5.057	1.00	37.15	A2	ATOM	674	CG	GIN	91	43.155	78.237	2.284	1.00	44.12	A2
ATOM	624	CB	TYR	86	42.598	70.255	-5.102	1.00	36.73	A2	ATOM	675	CD	GIN	91	44.348	78.769	1.542	1.00	46.96	A2
ATOM	625	CG	TYR	86	41.561	69.685	-6.081	1.00	33.66	A2	ATOM	676	OE1	GIN	91	45.235	78.083	1.068	1.00	47.42	A2
ATOM	626	CD1	TYR	86	41.946	69.312	-7.374	1.00	30.03	A2	ATOM	677	NE2	GIN	91	43.376	80.092	1.141	1.00	46.82	A2
ATOM	627	CE1	TYR	86	40.991	68.885	-8.280	1.00	30.08	A2	ATOM	678	HE2	GIN	91	43.690	80.685	1.700	1.00	0.00	A2
ATOM	628	CD2	TYR	86	40.224	69.623	-5.666	1.00	32.61	A2	ATOM	679	HE2	GIN	91	45.108	80.331	0.741	1.00	0.00	A2
ATOM	629	CE2	TYR	86	39.763	69.203	-6.574	1.00	31.66	A2	ATOM	680	C	GIN	91	40.129	77.231	2.061	1.00	37.22	A2
ATOM	630	CZ	TYR	86	39.656	68.818	-7.868	1.00	30.57	A2	ATOM	681	O	GIN	91	39.718	77.530	3.186	1.00	36.21	A2
ATOM	631	OH	TYR	86	38.670	68.428	-8.751	1.00	28.18	A2	ATOM	682	N	ALA	92	39.456	77.570	0.943	1.00	38.63	A2
ATOM	632	III	TYR	86	39.107	67.994	-9.485	1.00	0.00	A2	ATOM	683	II	ALA	92	39.808	77.205	0.098	1.00	0.00	A2
ATOM	633	C	TYR	86	43.054	72.318	-3.746	1.00	37.75	A2	ATOM	684	CA	ALA	92	38.243	78.402	0.880	1.00	36.10	A2
ATOM	634	O	TYR	86	42.173	72.469	-2.889	1.00	39.52	A2	ATOM	685	CB	ALA	92	37.657	78.436	-0.511	1.00	36.76	A2
ATOM	635	N	GIN	87	44.347	72.655	-3.478	1.00	36.93	A2	ATOM	686	C	ALA	92	37.139	77.905	1.770	1.00	38.95	A2
ATOM	636	II	GIN	87	45.044	72.463	-4.140	1.00	0.00	A2	ATOM	687	O	ALA	92	36.794	78.687	2.194	1.00	42.45	A2
ATOM	637	CA	GIN	87	44.749	73.332	-2.205	1.00	36.40	A2	ATOM	688	N	LEU	93	37.151	76.618	2.123	1.00	38.44	A2
ATOM	638	CB	GIN	87	46.210	73.668	-2.255	1.00	39.36	A2	ATOM	689	II	LEU	93	37.855	76.040	1.759	1.00	0.00	A2
ATOM	639	CG	GIN	87	47.126	72.993	-1.237	1.00	46.99	A2	ATOM	690	CA	LEU	93	36.111	76.018	2.972	1.00	36.90	A2
ATOM	640	CD	GIN	87	48.641	73.062	-1.576	1.00	50.96	A2	ATOM	691	CB	LEU	93	36.088	74.463	2.794	1.00	35.34	A2
ATOM	641	OE1	GIN	87	49.144	72.623	-2.627	1.00	52.15	A2	ATOM	692	CG	LEU	93	35.725	73.992	1.378	1.00	33.55	A2
ATOM	642	NE2	GIN	87	49.446	73.608	-0.663	1.00	52.96	A2	ATOM	693	CD1	LEU	93	36.159	72.583	1.129	1.00	33.26	A2
ATOM	643	HE2	GIN	87	49.055	73.957	0.164	1.00	0.00	A2	ATOM	694	CD2	LEU	93	34.254	74.167	1.215	1.00	32.16	A2
ATOM	644	HE2	GIN	87	50.396	73.621	-0.888	1.00	0.00	A2	ATOM	695	C	LEU	93	36.264	76.353	4.426	1.00	36.41	A2
ATOM	645	C	GIN	87	43.941	74.652	-2.013	1.00	34.36	A2	ATOM	696	O	LEU	93	35.473	75.917	5.256	1.00	35.17	A2
ATOM	646	O	GIN	87	43.414	74.990	-0.935	1.00	31.55	A2	ATOM	697	N	GLU	94	37.357	77.019	4.736	1.00	38.19	A2
ATOM	647	N	GLY	88	43.740	75.335	-3.159	1.00	32.73	A2	ATOM	698	II	GLU	94	38.022	77.167	4.035	1.00	0.00	A2
ATOM	648	II	GLY	88	44.165	75.005	-3.981	1.00	0.00	A2	ATOM	699	CA	GLU	94	37.627	77.573	6.038	1.00	42.71	A2
ATOM	649	CA	GLY	88	41.948	76.546	-3.232	1.00	30.81	A2	ATOM	700	CB	GLU	94	36.931	78.947	6.165	1.00	47.18	A2
ATOM	650	C	GLY	88	41.540	76.275	-2.731	1.00	30.47	A2	ATOM	701	CG	GLU	94	37.418	80.011	5.131	1.00	56.10	A2
ATOM	651	O	GLY	88	41.130	76.819	-1.703	1.00	30.27	A2	ATOM	702	CD	GLU	94	36.423	81.153	4.862	1.00	60.26	A2
ATOM	652	N	LEU	89	40.802	75.387	-3.406	1.00	29.01	A2	ATOM	703	OE1	GLU	94	35.728	81.109	3.023	1.00	60.76	A2
ATOM	653	II	LEU	89	41.220	74.912	-4.154	1.00	0.00	A2	ATOM	704	OE2	GLU	94	36.331	82.054	5.721	1.00	61.64	A2
ATOM	654	CA	LEU	89	39.447	75.102	-3.009	1.00	27.60	A2	ATOM	705	C	GLU	94	37.245	76.701	7.198	1.00	43.90	A2
ATOM	655	CB	LEU	89	38.922	74.073	-3.935	1.00	28.13	A2	ATOM	706	O	GLU	94	36.624	77.172	8.167	1.00	45.70	A2
ATOM	656	CG	LEU	89	38.764	74.583	-5.340	1.00	29.51	A2	ATOM	707	N	GLY	95	37.641	75.410	7.001	1.00	44.04	A2
ATOM	657	CD1	LEU	89	38.363	73.530	-6.364	1.00	24.13	A2	ATOM	708	II	GLY	95	38.024	75.192	6.127	1.00	0.00	A2
ATOM	658	CD2	LEU	89	37.673	75.637	-5.220	1.00	32.87	A2	ATOM	709	CA	GLY	95	37.519	74.310	7.981	1.00	42.49	A2
ATOM	659	C	LEU	89	39.352	74.629	-1.583	1.00	29.88	A2	ATOM	710	C	GLY	95	36.162	73.612	8.061	1.00	42.24	A2
ATOM	660	O	LEU	89	38.427	75.012	-0.860	1.00	30.81	A2	ATOM	711	O	GLY	95	36.028	72.596	8.739	1.00	40.02	A2
ATOM	661	N	LEU	90	40.317	73.839	-1.094	1.00	32.59	A2	ATOM	712	N	ILE	96	35.160	74.123	7.328	1.00	42.82	A2

FIGURE 5

ATOM	713	H	ILE	96	35.357	74.944	6.841	1.00	0.00	A2	ATOM	764	CG	PRO	102	40.799	68.687	13.776	1.00	41.02	A2
ATOM	714	CA	ILE	96	33.760	73.692	7.312	1.00	42.12	A2	ATOM	765	C	PRO	102	41.364	67.795	10.331	1.00	37.15	A2
ATOM	715	CB	ILE	96	33.665	72.233	6.800	1.00	36.33	A2	ATOM	766	O	PRO	102	42.358	67.854	9.600	1.00	38.88	A2
ATOM	716	CG2	ILE	96	32.248	71.768	6.789	1.00	34.79	A2	ATOM	767	N	THR	103	40.223	67.167	10.045	1.00	35.16	A2
ATOM	717	CG1	ILE	96	34.091	72.157	5.374	1.00	35.35	A2	ATOM	768	H	THR	103	39.466	67.223	10.662	1.00	0.00	A2
ATOM	718	CD	ILE	96	34.051	70.743	4.738	1.00	33.64	A2	ATOM	769	CA	THR	103	40.051	66.386	8.843	1.00	34.62	A2
ATOM	719	C	ILE	96	33.106	73.863	8.709	1.00	44.74	A2	ATOM	770	CB	THR	103	38.592	65.888	8.715	1.00	34.07	A2
ATOM	720	O	ILE	96	32.220	74.716	8.841	1.00	44.59	A2	ATOM	771	CG1	THR	103	38.011	65.896	10.548	1.00	0.00	A2
ATOM	721	N	SER	97	33.467	73.154	9.780	1.00	46.84	A2	ATOM	772	HG1	THR	103	38.312	64.896	7.594	1.00	31.79	A2
ATOM	722	H	SER	97	34.243	72.553	9.706	1.00	0.00	A2	ATOM	773	CG2	THR	103	40.417	67.215	7.425	1.00	34.61	A2
ATOM	723	CA	SER	97	32.900	73.359	11.105	1.00	48.91	A2	ATOM	774	C	THR	103	41.091	66.665	6.738	1.00	38.16	A2
ATOM	724	CB	SER	97	31.804	72.343	11.347	1.00	49.60	A2	ATOM	775	O	THR	103	40.054	68.498	7.529	1.00	32.49	A2
ATOM	725	CG	SER	97	32.211	71.120	11.954	1.00	52.85	A2	ATOM	776	N	LEU	104	39.504	68.923	8.229	1.00	0.00	A2
ATOM	726	HG	SER	97	31.406	70.573	11.942	1.00	0.00	A2	ATOM	777	H	LEU	104	40.471	69.267	6.370	1.00	30.49	A2
ATOM	727	C	SER	97	34.045	73.143	12.077	1.00	50.64	A2	ATOM	778	CA	LEU	104	39.616	70.430	6.242	1.00	31.51	A2
ATOM	728	O	SER	97	35.035	72.538	11.678	1.00	52.78	A2	ATOM	779	CB	LEU	104	38.356	69.996	5.611	1.00	36.61	A2
ATOM	729	N	PRO	98	34.063	73.474	13.348	1.00	52.12	A2	ATOM	780	CG	LEU	104	37.222	70.621	6.381	1.00	39.43	A2
ATOM	730	CD	PRO	98	33.002	74.170	14.016	1.00	52.90	A2	ATOM	781	CD1	LEU	104	38.418	70.294	4.132	1.00	37.89	A2
ATOM	731	CA	PRO	98	35.195	73.200	14.257	1.00	54.94	A2	ATOM	782	CD2	LEU	104	41.904	69.727	6.414	1.00	28.48	A2
ATOM	732	CB	PRO	98	34.750	73.717	15.600	1.00	54.78	A2	ATOM	783	C	LEU	104	42.583	69.825	5.398	1.00	28.47	A2
ATOM	733	CG	PRO	98	33.772	74.777	15.182	1.00	55.48	A2	ATOM	784	O	LEU	104	42.449	69.949	7.574	1.00	26.99	A2
ATOM	734	C	PRO	98	35.591	71.723	14.336	1.00	56.75	A2	ATOM	785	N	ASP	105	41.903	69.912	8.388	1.00	0.00	A2
ATOM	735	O	PRO	98	36.738	71.274	14.468	1.00	57.85	A2	ATOM	786	H	ASP	105	43.822	70.307	7.613	1.00	28.67	A2
ATOM	736	N	GLU	99	34.509	70.971	14.114	1.00	58.21	A2	ATOM	787	CA	ASP	105	44.139	70.584	9.038	1.00	33.06	A2
ATOM	737	H	GLU	99	33.652	71.400	14.078	1.00	0.00	A2	ATOM	788	CB	ASP	105	43.438	71.808	9.593	1.00	35.46	A2
ATOM	738	CA	GLU	99	34.543	69.537	14.281	1.00	58.48	A2	ATOM	789	CG	ASP	105	43.085	72.726	8.836	1.00	38.42	A2
ATOM	739	CB	GLU	99	33.111	69.104	14.304	1.00	63.30	A2	ATOM	790	CD1	ASP	105	43.244	71.816	10.808	1.00	39.10	A2
ATOM	740	CG	GLU	99	32.958	67.702	14.852	1.00	71.04	A2	ATOM	791	CD2	ASP	105	44.701	69.206	7.032	1.00	28.90	A2
ATOM	741	CD	GLU	99	32.076	66.838	13.962	1.00	76.95	A2	ATOM	792	C	ASP	105	45.551	69.479	6.175	1.00	29.62	A2
ATOM	742	OE1	GLU	99	32.209	65.608	14.079	1.00	80.63	A2	ATOM	793	O	ASP	105	44.415	67.950	7.401	1.00	26.86	A2
ATOM	743	OE2	GLU	99	31.295	67.382	13.153	1.00	77.99	A2	ATOM	794	N	THR	106	43.674	67.826	8.029	1.00	0.00	A2
ATOM	744	C	GLU	99	35.298	69.025	13.074	1.00	55.31	A2	ATOM	795	H	THR	106	45.143	66.770	6.935	1.00	24.81	A2
ATOM	745	O	GLU	99	36.251	68.270	13.210	1.00	55.96	A2	ATOM	796	CA	THR	106	44.558	65.456	7.477	1.00	26.03	A2
ATOM	746	N	LEU	100	34.916	69.475	11.891	1.00	51.23	A2	ATOM	797	CB	THR	106	44.680	65.566	8.894	1.00	31.53	A2
ATOM	747	H	LEU	100	34.214	70.159	11.841	1.00	0.00	A2	ATOM	798	CG1	THR	106	44.089	66.223	9.242	1.00	0.00	A2
ATOM	748	CA	LEU	100	35.577	69.052	10.678	1.00	48.08	A2	ATOM	799	CG2	THR	106	45.073	66.884	5.460	1.00	23.75	A2
ATOM	749	CB	LEU	100	34.627	69.341	9.574	1.00	45.52	A2	ATOM	800	CG1	THR	106	46.065	66.411	4.812	1.00	24.68	A2
ATOM	750	CG	LEU	100	33.544	68.337	9.674	1.00	45.39	A2	ATOM	801	C	THR	106	43.887	66.917	4.946	1.00	24.30	A2
ATOM	751	CD1	LEU	100	32.207	68.972	9.458	1.00	46.40	A2	ATOM	802	O	THR	106	43.145	67.176	5.528	1.00	0.00	A2
ATOM	752	CD2	LEU	100	33.851	67.245	8.677	1.00	47.48	A2	ATOM	803	N	LEU	107	43.668	66.783	3.531	1.00	27.29	A2
ATOM	753	C	LEU	100	36.956	69.629	10.368	1.00	46.77	A2	ATOM	804	H	LEU	107	42.158	66.913	3.273	1.00	25.45	A2
ATOM	754	O	LEU	100	37.578	69.244	9.357	1.00	46.62	A2	ATOM	805	CA	LEU	107	41.642	66.888	1.863	1.00	26.24	A2
ATOM	755	N	GLY	101	37.441	70.505	11.272	1.00	45.40	A2	ATOM	806	CB	LEU	107	42.095	65.649	1.158	1.00	26.41	A2
ATOM	756	H	GLY	101	36.893	70.704	12.056	1.00	0.00	A2	ATOM	807	CD1	LEU	107	40.140	66.925	1.914	1.00	27.42	A2
ATOM	757	CA	GLY	101	38.703	71.238	11.126	1.00	42.52	A2	ATOM	808	CD2	LEU	107	44.485	67.848	2.819	1.00	28.01	A2
ATOM	758	C	GLY	101	39.885	70.334	10.798	1.00	40.73	A2	ATOM	809	C	LEU	107	45.154	67.555	1.823	1.00	10.71	A2
ATOM	759	O	GLY	101	40.475	70.402	9.710	1.00	40.69	A2	ATOM	810	O	LEU	107	44.540	69.055	3.373	1.00	28.52	A2
ATOM	760	N	PRO	102	40.250	69.441	11.708	1.00	38.61	A2	ATOM	811	O	GLN	108	44.030	69.221	4.194	1.00	0.00	A2
ATOM	761	CD	PRO	102	39.676	69.350	13.027	1.00	39.76	A2	ATOM	812	N	GLN	108	45.343	70.132	2.792	1.00	28.38	A2
ATOM	762	CA	PRO	102	41.390	68.566	11.606	1.00	37.30	A2											
ATOM	763	CB	PRO	102	41.294	67.690	12.775	1.00	39.36	A2											

ATOM	815	CB	GIN	108	45.138	71.363	3.630	1.00	30.15	A2	ATOM	866	II	PIIE	114	51.344	66.164	-1.361	1.00	0.00	A2
ATOM	816	CG	GIN	108	43.711	71.787	3.542	1.00	32.27	A2	ATOM	867	CA	PIIE	114	52.109	65.328	-3.103	1.00	27.84	A2
ATOM	817	CG	GIN	108	43.606	73.192	4.048	1.00	35.64	A2	ATOM	868	CB	PIIE	114	50.708	64.794	-3.226	1.00	21.18	A2
ATOM	818	OEL	GIN	108	43.085	73.484	5.175	1.00	36.07	A2	ATOM	869	CG	PIIE	114	50.565	63.928	-4.420	1.00	21.04	A2
ATOM	819	NE2	GIN	108	44.189	74.044	3.213	1.00	33.58	A2	ATOM	870	CDI	PIIE	114	51.623	63.225	-4.938	1.00	24.05	A2
ATOM	820	IEE1	GIN	108	44.582	73.701	2.386	1.00	0.00	A2	ATOM	871	CDI	PIIE	114	49.369	63.914	-5.046	1.00	22.37	A2
ATOM	821	IEE2	GIN	108	44.895	74.986	3.471	1.00	0.00	A2	ATOM	872	CE2	PIIE	114	51.476	62.514	-6.102	1.00	24.54	A2
ATOM	822	C	GIN	108	46.840	69.842	2.675	1.00	26.40	A2	ATOM	873	CE2	PIIE	114	49.211	63.207	-6.212	1.00	21.15	A2
ATOM	823	N	GIN	108	47.450	69.955	1.597	1.00	27.57	A2	ATOM	874	CZ	PIIE	114	50.263	62.509	-6.741	1.00	24.71	A2
ATOM	824	N	GIN	108	47.388	69.473	3.833	1.00	25.81	A2	ATOM	875	C	PIIE	114	52.453	66.291	-4.190	1.00	29.20	A2
ATOM	825	II	LEU	109	46.795	69.495	4.615	1.00	0.00	A2	ATOM	876	C	PIIE	114	53.072	65.883	-5.158	1.00	40.84	A2
ATOM	826	CA	LEU	109	48.764	69.003	4.043	1.00	27.96	A2	ATOM	877	N	ALIA	115	52.057	67.554	-4.058	1.00	31.99	A2
ATOM	827	CB	LEU	109	48.951	68.637	5.513	1.00	29.41	A2	ATOM	878	II	ALIA	115	51.446	67.768	-3.317	1.00	0.00	A2
ATOM	828	CG	LEU	109	48.712	69.171	6.520	1.00	31.78	A2	ATOM	879	CA	ALIA	115	52.423	68.655	-4.952	1.00	31.29	A2
ATOM	829	CD1	LEU	109	48.750	69.688	7.933	1.00	29.16	A2	ATOM	880	CB	ALIA	115	51.824	69.939	-4.420	1.00	30.65	A2
ATOM	830	CD2	LEU	109	49.724	70.889	6.285	1.00	31.19	A2	ATOM	881	C	ALIA	115	53.936	68.787	-4.976	1.00	31.31	A2
ATOM	831	C	LEU	109	49.168	67.790	3.186	1.00	26.80	A2	ATOM	882	O	ALIA	115	54.539	68.873	-6.044	1.00	40.36	A2
ATOM	832	O	LEU	109	50.214	67.791	2.544	1.00	26.81	A2	ATOM	883	N	THIR	116	54.551	68.846	-3.813	1.00	32.20	A2
ATOM	833	N	ASP	110	48.305	66.807	3.090	1.00	25.98	A2	ATOM	884	II	THIR	116	54.013	68.910	-2.992	1.00	0.00	A2
ATOM	834	N	ASP	110	47.471	66.835	3.600	1.00	0.00	A2	ATOM	885	CA	THIR	116	55.998	68.997	-3.656	1.00	34.91	A2
ATOM	835	CA	ASP	110	48.590	65.684	2.250	1.00	23.32	A2	ATOM	886	CB	THIR	116	56.325	68.953	-2.150	1.00	35.78	A2
ATOM	836	CB	ASP	110	47.577	64.570	2.553	1.00	26.34	A2	ATOM	887	CG1	THIR	116	55.564	70.038	-1.576	1.00	35.58	A2
ATOM	837	CG	ASP	110	47.905	63.878	3.894	1.00	31.10	A2	ATOM	888	HG1	THIR	116	54.942	69.844	-0.939	1.00	0.00	A2
ATOM	838	CD1	ASP	110	47.070	63.093	4.323	1.00	34.98	A2	ATOM	889	CG2	THIR	116	57.816	67.726	-1.921	1.00	35.38	A2
ATOM	839	CD2	ASP	110	48.958	64.107	4.555	1.00	34.06	A2	ATOM	890	C	THIR	116	56.714	67.726	-4.304	1.00	37.14	A2
ATOM	840	C	ASP	110	48.557	66.138	0.842	1.00	20.31	A2	ATOM	891	O	THIR	116	57.641	67.937	-5.066	1.00	39.27	A2
ATOM	841	O	ASP	110	49.493	65.711	0.165	1.00	20.63	A2	ATOM	892	N	THIR	117	56.318	66.485	-4.045	1.00	49.05	A2
ATOM	842	N	VAL	111	47.627	66.998	0.363	1.00	20.80	A2	ATOM	893	II	THIR	117	55.615	66.383	-3.369	1.00	0.00	A2
ATOM	843	N	VAL	111	46.900	67.310	0.944	1.00	0.00	A2	ATOM	894	CA	THIR	117	56.840	65.269	-4.630	1.00	40.23	A2
ATOM	844	CA	VAL	111	47.771	67.454	-1.019	1.00	20.44	A2	ATOM	895	CB	THIR	117	55.909	64.920	-4.216	1.00	49.99	A2
ATOM	845	CB	VAL	111	46.531	68.364	-1.376	1.00	23.60	A2	ATOM	896	CG1	THIR	117	56.149	63.920	-2.820	1.00	41.66	A2
ATOM	846	CG1	VAL	111	45.668	68.946	-2.808	1.00	23.04	A2	ATOM	897	HG1	THIR	117	55.653	64.559	-2.286	1.00	0.00	A2
ATOM	847	CG2	VAL	111	45.289	67.497	-1.371	1.00	24.30	A2	ATOM	898	CG2	THIR	117	56.110	62.781	-4.981	1.00	38.23	A2
ATOM	848	C	VAL	111	49.006	68.224	-1.245	1.00	20.82	A2	ATOM	899	C	THIR	117	56.882	65.417	-6.134	1.00	43.23	A2
ATOM	849	O	VAL	111	49.617	68.006	-2.303	1.00	19.22	A2	ATOM	900	O	THIR	117	57.934	65.253	-6.749	1.00	46.29	A2
ATOM	850	N	ALA	112	49.442	69.063	-0.267	1.00	21.84	A2	ATOM	901	N	ILE	118	55.763	63.777	-6.741	1.00	45.87	A2
ATOM	851	H	ALA	112	48.839	69.590	0.492	1.00	0.00	A2	ATOM	902	II	ILE	118	54.962	65.942	-6.200	1.00	0.00	A2
ATOM	852	CA	ALA	112	50.708	69.805	-0.295	1.00	24.16	A2	ATOM	903	CA	ILE	118	53.659	65.914	-8.182	1.00	47.97	A2
ATOM	853	CB	ALA	112	50.861	70.561	1.011	1.00	22.69	A2	ATOM	904	CB	ILE	118	54.170	66.271	-8.452	1.00	47.69	A2
ATOM	854	C	ALA	112	51.931	68.878	-0.486	1.00	28.58	A2	ATOM	905	CG2	ILE	118	54.041	66.930	-9.835	1.00	47.03	A2
ATOM	855	O	ALA	112	52.778	69.026	-1.390	1.00	32.53	A2	ATOM	906	CG1	ILE	118	53.302	65.011	-8.244	1.00	44.03	A2
ATOM	856	N	ASP	113	52.086	67.852	0.343	1.00	30.21	A2	ATOM	907	CD	ILE	118	53.651	63.883	-9.236	1.00	41.71	A2
ATOM	857	CA	ASP	113	53.084	66.847	1.130	1.00	0.00	A2	ATOM	908	C	ILE	118	56.647	66.932	-8.774	1.00	50.69	A2
ATOM	858	CB	ASP	113	53.084	66.946	0.166	1.00	31.70	A2	ATOM	909	O	ILE	118	57.390	66.676	-9.681	1.00	49.98	A2
ATOM	859	CB	ASP	113	52.706	65.659	0.953	1.00	36.31	A2	ATOM	910	N	TRP	119	56.697	68.061	-8.015	1.00	54.68	A2
ATOM	860	CG	ASP	113	53.170	65.758	2.357	1.00	42.27	A2	ATOM	911	II	TRP	119	56.164	68.135	-7.197	1.00	0.00	A2
ATOM	861	OD1	ASP	113	52.559	65.109	3.289	1.00	46.37	A2	ATOM	912	CA	TRP	119	57.575	69.142	-7.477	1.00	58.98	A2
ATOM	862	OD2	ASP	113	54.160	66.461	2.589	1.00	48.93	A2	ATOM	913	CB	TRP	119	57.302	70.367	-7.977	1.00	59.84	A2
ATOM	863	C	ASP	113	53.315	66.361	-1.239	1.00	32.82	A2	ATOM	914	CG	TRP	119	58.051	71.579	-8.196	1.00	62.64	A2
ATOM	864	O	ASP	113	54.433	66.308	-1.734	1.00	36.25	A2	ATOM	915	CD2	TRP	119	57.596	72.721	-9.307	1.00	64.78	A2
ATOM	865	N	ASP	114	52.187	65.978	-1.850	1.00	30.24	A2	ATOM	916	CD2	TRP	119	58.699	72.955	-9.643	1.00	62.55	A2

FIGURE 5

ATOM	917	CE3 TRP	119	56.465	72.314	-10.080	1.00	66.02	ATOM	968	O GLU	123	65.092	68.558	-12.356	1.00	78.22	A2
ATOM	918	CD1 TRP	119	59.322	71.870	-7.863	1.00	64.12	ATOM	969	N GLU	124	64.504	67.110	-10.765	1.00	77.66	A2
ATOM	919	NE1 TRP	119	59.680	72.727	-8.784	1.00	65.00	ATOM	970	II GLU	124	63.867	66.852	-10.060	1.00	0.00	A2
ATOM	920	HE1 TRP	119	60.568	73.140	-8.828	1.00	0.00	ATOM	971	CA GLU	124	65.574	66.215	-11.167	1.00	78.37	A2
ATOM	921	C22 TRP	119	58.726	73.794	-10.714	1.00	62.90	ATOM	972	CB GLU	124	65.600	65.051	-10.195	1.00	80.79	A2
ATOM	922	C23 TRP	119	56.469	73.157	-11.170	1.00	65.18	ATOM	973	CG GLU	124	64.387	64.132	-10.150	1.00	81.29	A2
ATOM	923	CH2 TRP	119	57.591	73.887	-11.481	1.00	64.40	ATOM	974	CD GLU	124	64.375	63.248	-8.908	1.00	85.51	A2
ATOM	924	C TRP	119	59.021	68.664	-8.352	1.00	61.26	ATOM	975	OEL GLU	124	64.733	63.729	-7.824	1.00	80.44	A2
ATOM	925	O TRP	119	59.748	68.788	-9.343	1.00	62.12	ATOM	976	OEL GLU	124	64.006	62.075	-9.024	1.00	86.59	A2
ATOM	926	N GLN	120	59.447	68.065	-7.249	1.00	62.91	ATOM	977	C GLU	124	65.534	65.705	-12.612	1.00	78.01	A2
ATOM	927	H GLN	120	58.811	67.961	-6.519	1.00	0.00	ATOM	978	O GLU	124	66.480	65.057	-13.060	1.00	78.91	A2
ATOM	928	CA GLN	120	60.786	67.504	-7.113	1.00	65.16	ATOM	979	N LEU	125	64.460	65.943	-13.363	1.00	77.11	A2
ATOM	929	CB GLN	120	60.900	66.800	-5.780	1.00	66.56	ATOM	980	II LEU	125	63.666	66.340	-12.945	1.00	0.00	A2
ATOM	930	CG GLN	120	60.627	67.678	-4.582	1.00	67.18	ATOM	981	CA LEU	125	64.387	65.583	-14.771	1.00	76.23	A2
ATOM	931	CD GLN	120	60.725	66.907	-3.284	1.00	67.77	ATOM	982	CB LEU	125	63.061	64.832	-14.952	1.00	76.88	A2
ATOM	932	OEL GLN	120	61.221	67.465	-2.319	1.00	69.31	ATOM	983	CG LEU	125	62.392	64.382	-16.263	1.00	76.63	A2
ATOM	933	NE2 GLN	120	60.305	65.654	-3.129	1.00	67.39	ATOM	984	CD1 LEU	125	63.350	63.754	-17.276	1.00	76.67	A2
ATOM	934	HE21 GLN	120	59.903	65.174	-3.877	1.00	0.00	ATOM	985	CD2 LEU	125	61.309	63.402	-15.839	1.00	75.89	A2
ATOM	935	HE22 GLN	120	60.441	65.282	-2.234	1.00	0.00	ATOM	986	C LEU	125	64.506	66.827	-15.648	1.00	75.84	A2
ATOM	936	C GLN	120	61.169	66.509	-8.722	1.00	66.22	ATOM	987	O LEU	125	64.360	66.788	-16.871	1.00	75.36	A2
ATOM	937	O GLN	120	62.326	66.421	-8.662	1.00	66.50	ATOM	988	N GLY	126	64.759	67.968	-15.027	1.00	75.90	A2
ATOM	938	N GLN	121	60.202	65.745	-8.706	1.00	67.10	ATOM	989	II GLY	126	64.741	67.976	-14.056	1.00	0.00	A2
ATOM	939	H GLN	121	59.307	65.754	-8.303	1.00	0.00	ATOM	990	CA GLY	126	64.968	69.213	-15.736	1.00	77.58	A2
ATOM	940	CA GLN	121	60.480	64.878	-9.812	1.00	68.66	ATOM	991	C GLY	126	63.677	69.814	-16.330	1.00	78.63	A2
ATOM	941	CB GLN	121	59.292	63.971	-10.070	1.00	67.96	ATOM	992	O GLY	126	63.735	70.736	-17.146	1.00	78.55	A2
ATOM	942	CG GLN	121	59.614	62.937	-11.128	1.00	68.89	ATOM	993	N MET	127	62.524	69.343	-15.933	1.00	81.08	A2
ATOM	943	CD GLN	121	60.940	62.236	-10.852	1.00	71.37	ATOM	994	II MET	127	62.522	68.603	-15.293	1.00	0.00	A2
ATOM	944	OEL GLN	121	61.212	61.706	-9.777	1.00	71.70	ATOM	995	CA MET	127	61.266	69.902	-16.415	1.00	81.46	A2
ATOM	945	NE2 GLN	121	61.879	62.262	-11.786	1.00	74.41	ATOM	996	CG MET	127	60.191	68.802	-16.361	1.00	81.86	A2
ATOM	946	HE21 GLN	121	61.707	62.729	-12.627	1.00	0.00	ATOM	997	CG MET	127	60.708	67.599	-17.147	1.00	82.66	A2
ATOM	947	HE22 GLN	121	62.736	61.859	-11.541	1.00	0.00	ATOM	998	SD MET	127	59.682	66.115	-17.282	1.00	81.70	A2
ATOM	948	C GLN	121	60.760	65.743	-11.045	1.00	70.48	ATOM	999	CE MET	127	60.236	65.620	-18.900	1.00	81.23	A2
ATOM	949	O GLN	121	61.671	65.436	-11.827	1.00	70.94	ATOM	1000	C MET	127	60.847	71.131	-15.599	1.00	82.18	A2
ATOM	950	N MET	122	60.019	66.846	-11.236	1.00	71.67	ATOM	1001	OT1 MET	127	60.116	71.958	-16.142	1.00	83.86	A2
ATOM	951	H MET	122	59.351	67.087	-10.535	1.00	0.00	ATOM	1002	OT2 MET	127	61.267	71.285	-14.446	1.00	82.04	A2
ATOM	952	CA MET	122	60.190	67.688	-12.412	1.00	72.62	ATOM	1003	CB MET	138	39.323	80.595	-4.492	1.00	50.39	A1
ATOM	953	CB MET	122	59.123	68.818	-12.448	1.00	73.12	ATOM	1004	CG MET	138	40.123	79.298	-4.421	1.00	51.97	A1
ATOM	954	CG MET	122	57.880	68.343	-13.083	1.00	73.64	ATOM	1005	SD MET	138	40.561	78.973	-6.145	1.00	60.85	A1
ATOM	955	SD MET	122	56.669	69.662	-13.295	1.00	75.44	ATOM	1006	CE MET	138	41.129	77.310	-6.351	1.00	61.48	A1
ATOM	956	CE MET	122	55.695	69.349	-11.861	1.00	76.43	ATOM	1007	C MET	138	37.021	81.072	-5.454	1.00	60.26	A1
ATOM	957	C MET	122	61.566	68.281	-12.411	1.00	73.22	ATOM	1008	O MET	138	36.832	82.262	-5.181	1.00	62.98	A1
ATOM	958	O MET	122	62.240	68.287	-13.441	1.00	73.03	ATOM	1009	HT1 MET	138	38.497	82.600	-6.075	1.00	0.00	A1
ATOM	959	N GLU	123	61.991	68.697	-11.223	1.00	74.74	ATOM	1010	HT2 MET	138	38.313	81.757	-7.529	1.00	0.00	A1
ATOM	960	H GLU	123	61.372	68.617	-10.466	1.00	0.00	ATOM	1011	N MET	138	38.839	81.784	-6.639	1.00	60.49	A1
ATOM	961	CA GLU	123	63.305	69.262	-11.018	1.00	75.95	ATOM	1012	HT3 MET	138	39.865	81.816	-6.768	1.00	0.00	A1
ATOM	962	CB GLU	123	63.484	69.665	-9.597	1.00	75.72	ATOM	1013	CA MET	138	38.445	80.672	-5.787	1.00	60.51	A1
ATOM	963	CG GLU	123	62.644	70.906	-9.500	1.00	79.11	ATOM	1014	N PKO	139	35.995	80.242	-5.612	1.00	57.82	A1
ATOM	964	CD GLU	123	62.651	71.529	-8.122	1.00	83.02	ATOM	1015	CD PKO	139	36.028	79.060	-6.448	1.00	58.10	A1
ATOM	965	OEL GLU	123	62.741	72.763	-8.057	1.00	84.15	ATOM	1016	CA PKO	139	34.654	80.538	-5.142	1.00	54.67	A1
ATOM	966	HE2 GLU	123	62.543	70.789	-7.133	1.00	84.45	ATOM	1017	CB PKO	139	34.870	79.323	-5.525	1.00	54.54	A1
ATOM	967	C GLU	123	64.381	68.280	-11.386	1.00	77.17	ATOM	1018	CG PKO	139	34.945	78.290	-5.755	1.00	58.20	A1

FIGURE 5

ATOM 1019 C PRO 139	34.588	80.875	-3.664	1.00	52.24	A3
ATOM 1020 O PRO 139	35.507	80.623	-2.882	1.00	51.89	A3
ATOM 1021 N ALA 140	33.499	81.547	-3.342	1.00	49.86	A3
ATOM 1022 H ALA 140	32.789	81.676	-4.005	1.00	0.00	A3
ATOM 1023 CA ALA 140	33.234	81.926	-1.994	1.00	47.39	A3
ATOM 1024 CB ALA 140	32.966	83.413	-1.895	1.00	49.34	A3
ATOM 1025 C ALA 140	31.978	81.153	-1.590	1.00	49.25	A3
ATOM 1026 O ALA 140	30.889	81.162	-2.205	1.00	49.06	A3
ATOM 1027 N PHE 141	32.293	80.442	-0.506	1.00	47.48	A3
ATOM 1028 H PHE 141	33.190	80.550	-0.122	1.00	0.00	A3
ATOM 1029 CA PHE 141	31.401	79.552	0.208	1.00	45.66	A3
ATOM 1030 CB PHE 141	32.215	78.305	0.792	1.00	40.28	A3
ATOM 1031 CG PHE 141	32.684	77.404	-0.349	1.00	35.35	A3
ATOM 1032 CD1 PHE 141	31.800	76.591	-1.006	1.00	34.39	A3
ATOM 1033 CD2 PHE 141	33.966	77.497	-0.830	1.00	37.69	A3
ATOM 1034 CE1 PHE 141	32.174	75.895	-2.133	1.00	34.00	A3
ATOM 1035 CE2 PHE 141	34.358	76.807	-1.956	1.00	36.69	A3
ATOM 1036 CZ PHE 141	33.449	76.001	-2.614	1.00	37.29	A3
ATOM 1037 C PHE 141	31.003	80.580	1.242	1.00	46.54	A3
ATOM 1038 O PHE 141	31.584	80.664	2.317	1.00	48.26	A3
ATOM 1039 N ALA 142	30.067	81.452	0.843	1.00	47.38	A3
ATOM 1040 H ALA 142	29.624	81.295	-0.020	1.00	0.00	A3
ATOM 1041 CA ALA 142	29.581	82.564	1.668	1.00	46.06	A3
ATOM 1042 CB ALA 142	28.731	83.546	0.879	1.00	45.04	A3
ATOM 1043 C ALA 142	28.703	82.132	2.802	1.00	45.27	A3
ATOM 1044 O ALA 142	28.343	83.002	3.584	1.00	47.38	A3
ATOM 1045 N SER 143	28.318	80.860	2.899	1.00	43.36	A3
ATOM 1046 H SER 143	28.724	80.201	2.503	1.00	0.00	A3
ATOM 1047 CA SER 143	27.377	80.392	3.897	1.00	41.94	A3
ATOM 1048 CB SER 143	26.036	80.129	3.181	1.00	44.17	A3
ATOM 1049 CG SER 143	25.323	78.918	3.536	1.00	48.18	A3
ATOM 1050 HG SER 143	24.455	78.974	3.098	1.00	0.00	A3
ATOM 1051 C SER 143	27.877	79.145	4.602	1.00	39.79	A3
ATOM 1052 O SER 143	28.763	78.452	4.132	1.00	38.50	A3
ATOM 1053 N ALA 144	27.218	78.775	5.683	1.00	39.10	A3
ATOM 1054 H ALA 144	26.449	79.312	5.960	1.00	0.00	A3
ATOM 1055 CA ALA 144	27.566	77.586	6.411	1.00	39.22	A3
ATOM 1056 CB ALA 144	26.964	76.420	5.627	1.00	36.97	A3
ATOM 1057 C ALA 144	26.964	76.420	5.627	1.00	41.58	A3
ATOM 1058 O ALA 144	27.706	75.448	5.444	1.00	42.07	A3
ATOM 1059 N PHE 145	25.719	76.407	5.076	1.00	40.77	A3
ATOM 1060 H PHE 145	25.149	77.203	5.110	1.00	0.00	A3
ATOM 1061 CA PHE 145	25.307	75.234	4.312	1.00	39.31	A3
ATOM 1062 CB PHE 145	23.877	75.396	3.798	1.00	36.46	A3
ATOM 1063 CG PHE 145	23.477	74.452	2.641	1.00	31.91	A3
ATOM 1064 CD1 PHE 145	23.579	74.900	1.323	1.00	29.02	A3
ATOM 1065 CD2 PHE 145	23.013	73.185	2.916	1.00	29.40	A3
ATOM 1066 CE1 PHE 145	23.225	74.100	0.277	1.00	28.34	A3
ATOM 1067 CE2 PHE 145	22.661	72.389	1.858	1.00	28.80	A3
ATOM 1068 CZ PHE 145	22.764	72.831	0.549	1.00	30.58	A3
ATOM 1069 C PHE 145	26.266	75.071	3.120	1.00	40.44	A3
ATOM 1070 O PHE 145	26.556	73.938	2.697	1.00	40.55	A3
ATOM 1071 N GLN 146	26.745	76.232	2.619	1.00	49.11	A3
ATOM 1072 H GLN 146	26.437	77.073	3.015	1.00	0.00	A3
ATOM 1073 CA GLN 146	27.660	76.263	1.531	1.00	48.01	A3
ATOM 1074 CB GLN 146	27.807	77.644	1.054	1.00	48.92	A3
ATOM 1075 CG GLN 146	26.884	78.046	0.949	1.00	43.26	A3
ATOM 1076 CD GLN 146	27.171	79.440	-0.522	1.00	45.47	A3
ATOM 1077 CE1 GLN 146	27.851	80.253	0.043	1.00	47.57	A3
ATOM 1078 CE2 GLN 146	26.689	79.793	-1.092	1.00	47.50	A3
ATOM 1079 HE2 GLN 146	26.149	79.149	-2.190	1.00	0.00	A3
ATOM 1080 HE22 GLN 146	26.913	80.690	-2.021	1.00	0.00	A3
ATOM 1081 C GLN 146	29.005	75.670	1.836	1.00	47.45	A3
ATOM 1082 O GLN 146	29.634	75.093	0.950	1.00	48.28	A3
ATOM 1083 N ARG 147	29.511	75.775	3.054	1.00	46.37	A3
ATOM 1084 H ARG 147	29.044	76.300	3.738	1.00	0.00	A3
ATOM 1085 CA ARG 147	30.798	75.180	3.357	1.00	45.68	A3
ATOM 1086 CB ARG 147	31.799	75.574	4.713	1.00	47.12	A3
ATOM 1087 CG ARG 147	31.730	77.016	4.697	1.00	42.68	A3
ATOM 1088 CD ARG 147	32.034	77.494	6.093	1.00	49.54	A3
ATOM 1089 NE ARG 147	32.674	78.774	5.877	1.00	58.21	A3
ATOM 1090 HE ARG 147	32.475	79.252	5.045	1.00	0.00	A3
ATOM 1091 CZ ARG 147	33.519	79.373	6.742	1.00	62.77	A3
ATOM 1092 NH1 ARG 147	33.905	78.868	7.936	1.00	64.96	A3
ATOM 1093 NH11 ARG 147	34.545	79.379	8.510	1.00	0.00	A3
ATOM 1094 NH12 ARG 147	33.561	77.980	8.239	1.00	0.00	A3
ATOM 1095 NH2 ARG 147	31.960	80.584	6.403	1.00	64.80	A3
ATOM 1096 NH21 ARG 147	34.599	81.069	6.999	1.00	0.00	A3
ATOM 1097 NH22 ARG 147	33.665	80.996	5.541	1.00	0.00	A3
ATOM 1098 C ARG 147	30.570	73.707	3.337	1.00	44.91	A3
ATOM 1099 O ARG 147	31.233	73.050	2.539	1.00	44.56	A3
ATOM 1100 N ALA 148	29.544	73.194	4.040	1.00	44.44	A3
ATOM 1101 H ALA 148	28.926	73.818	4.482	1.00	0.00	A3
ATOM 1102 CA ALA 148	29.358	71.754	4.172	1.00	45.92	A3
ATOM 1103 CB ALA 148	28.217	71.426	5.153	1.00	52.85	A3
ATOM 1104 C ALA 148	29.077	71.095	2.845	1.00	44.40	A3
ATOM 1105 O ALA 148	29.765	70.141	2.157	1.00	44.31	A3
ATOM 1106 N ALA 149	28.169	71.657	2.077	1.00	42.60	A3
ATOM 1107 H ALA 149	27.662	72.424	2.411	1.00	0.00	A3
ATOM 1108 CA ALA 149	27.890	71.134	0.757	1.00	42.70	A3
ATOM 1109 CB ALA 149	26.595	71.774	0.299	1.00	41.51	A3
ATOM 1110 C ALA 149	29.032	71.381	-0.258	1.00	44.75	A3
ATOM 1111 O ALA 149	29.208	70.661	-1.264	1.00	44.49	A3
ATOM 1112 N GLY 150	29.867	72.401	-0.052	1.00	44.58	A3
ATOM 1113 H GLY 150	29.724	73.035	0.682	1.00	0.00	A3
ATOM 1114 CA GLY 150	31.017	72.608	-0.913	1.00	41.79	A3
ATOM 1115 C GLY 150	32.113	71.629	-0.478	1.00	41.99	A3
ATOM 1116 O GLY 150	32.997	71.261	-1.265	1.00	41.77	A3
ATOM 1117 N GLY 151	32.075	71.161	0.773	1.00	49.64	A3
ATOM 1118 H GLY 151	31.412	71.524	1.394	1.00	0.00	A3
ATOM 1119 CA GLY 151	33.018	70.166	1.243	1.00	42.16	A3
ATOM 1120 C GLY 151	32.764	68.969	0.409	1.00	44.98	A3

FIGURE 5

ATOM	1121	O	GLY	151	33.664	68.501	-0.349	1.00	35.66	A3	ATOM	1172	C	11S	157	37.291	65.476	-7.260	1.00	29.68	A3
ATOM	1122	N	VAL	152	31.486	68.418	0.451	1.00	31.87	A3	ATOM	1173	O	11S	157	37.950	65.059	-8.219	1.00	29.65	A3
ATOM	1123	H	VAL	152	30.867	68.906	1.040	1.00	0.00	A3	ATOM	1174	N	LEU	158	37.801	65.069	-6.071	1.00	29.24	A3
ATOM	1124	CA	VAL	152	30.978	67.240	-0.275	1.00	29.61	A3	ATOM	1175	H	LEU	158	37.213	65.901	-5.326	1.00	0.00	A3
ATOM	1125	CB	VAL	152	29.419	67.145	-0.125	1.00	27.63	A3	ATOM	1176	CA	LEU	158	39.216	65.479	-5.826	1.00	31.94	A3
ATOM	1126	CG1	VAL	152	28.083	66.035	-0.976	1.00	27.37	A3	ATOM	1177	CB	LEU	158	39.609	65.949	-4.373	1.00	28.66	A3
ATOM	1127	CG2	VAL	152	29.002	66.786	1.279	1.00	24.74	A3	ATOM	1178	CG	LEU	158	41.008	65.751	-3.859	1.00	24.32	A3
ATOM	1128	C	VAL	152	31.351	67.294	-1.762	1.00	29.91	A3	ATOM	1179	CD1	LEU	158	41.990	66.378	-4.776	1.00	20.87	A3
ATOM	1129	O	VAL	152	31.805	66.329	-2.393	1.00	31.75	A3	ATOM	1180	CD2	LEU	158	41.099	66.330	-2.477	1.00	24.86	A3
ATOM	1130	N	LEU	153	31.236	68.452	-2.361	1.00	29.26	A3	ATOM	1181	C	LEU	158	39.468	63.994	-6.027	1.00	31.46	A3
ATOM	1131	H	LEU	153	30.881	69.219	-1.860	1.00	0.00	A3	ATOM	1182	O	LEU	158	40.298	63.609	-6.844	1.00	30.58	A3
ATOM	1132	CA	LEU	153	31.559	68.607	-3.756	1.00	26.77	A3	ATOM	1183	N	GLN	159	38.652	63.225	-5.340	1.00	33.54	A3
ATOM	1133	CB	LEU	153	30.881	69.858	-4.160	1.00	28.22	A3	ATOM	1184	H	GLN	159	38.011	63.676	-4.748	1.00	0.00	A3
ATOM	1134	CG	LEU	153	29.943	69.894	-5.316	1.00	30.67	A3	ATOM	1185	CA	GLN	159	38.594	61.792	-5.447	1.00	35.73	A3
ATOM	1135	CD1	LEU	153	28.580	69.281	-5.090	1.00	26.48	A3	ATOM	1186	CB	GLN	159	37.308	61.492	-4.813	1.00	37.26	A3
ATOM	1136	CD2	LEU	153	29.741	71.365	-5.496	1.00	34.46	A3	ATOM	1187	CG	GLN	159	37.064	60.063	-4.520	1.00	45.01	A3
ATOM	1137	C	LEU	153	33.032	68.628	-4.111	1.00	26.08	A3	ATOM	1188	CD	GLN	159	37.755	59.611	-3.256	1.00	46.24	A3
ATOM	1138	O	LEU	153	33.419	68.187	-5.212	1.00	26.78	A3	ATOM	1189	OE1	GLN	159	38.142	58.443	-3.232	1.00	48.29	A3
ATOM	1139	N	VAL	154	33.902	69.180	-3.269	1.00	26.12	A3	ATOM	1190	NE2	GLN	159	37.936	60.456	-2.224	1.00	47.82	A3
ATOM	1140	H	VAL	154	33.589	69.557	-2.416	1.00	0.00	A3	ATOM	1191	HE21	GLN	159	37.575	61.364	-2.289	1.00	0.00	A3
ATOM	1141	CA	VAL	154	35.330	69.259	-3.611	1.00	26.23	A3	ATOM	1192	HE22	GLN	159	38.412	60.101	-1.447	1.00	0.00	A3
ATOM	1142	CB	VAL	154	36.057	70.299	-2.692	1.00	26.51	A3	ATOM	1193	C	GLN	159	38.686	61.381	-6.921	1.00	36.24	A3
ATOM	1143	CG1	VAL	154	37.578	70.188	-2.942	1.00	25.01	A3	ATOM	1194	O	GLN	159	39.632	60.690	-7.324	1.00	38.97	A3
ATOM	1144	CG2	VAL	154	35.528	71.728	-2.945	1.00	27.82	A3	ATOM	1195	N	SER	160	37.824	61.896	-7.796	1.00	35.48	A3
ATOM	1145	C	VAL	154	35.933	67.850	-3.375	1.00	26.80	A3	ATOM	1196	H	SER	160	37.142	62.540	-7.498	1.00	0.00	A3
ATOM	1146	O	VAL	154	36.678	67.363	-4.229	1.00	26.27	A3	ATOM	1197	CA	SER	160	37.869	61.564	-9.203	1.00	34.96	A3
ATOM	1147	N	ALA	155	35.635	67.241	-2.199	1.00	24.76	A3	ATOM	1198	CB	SER	160	36.645	62.100	-8.863	1.00	37.54	A3
ATOM	1148	H	ALA	155	35.084	67.758	-1.570	1.00	0.00	A3	ATOM	1199	CG	SER	160	35.587	62.434	-8.942	1.00	44.81	A3
ATOM	1149	CA	ALA	155	36.095	65.940	-1.782	1.00	25.21	A3	ATOM	1200	IG	SER	160	35.340	61.689	-8.387	1.00	0.00	A3
ATOM	1150	CB	ALA	155	35.463	65.572	-0.457	1.00	25.25	A3	ATOM	1201	C	SER	160	39.090	62.095	-9.922	1.00	31.65	A3
ATOM	1151	C	ALA	155	35.708	64.946	-2.841	1.00	26.94	A3	ATOM	1202	O	SER	160	39.605	61.382	-10.785	1.00	35.42	A3
ATOM	1152	O	ALA	155	36.594	64.288	-3.398	1.00	29.96	A3	ATOM	1203	N	PIE	161	39.615	63.293	-9.595	1.00	33.82	A3
ATOM	1153	N	SER	156	34.450	64.982	-3.282	1.00	29.96	A3	ATOM	1204	H	PIE	161	39.203	63.796	-8.864	1.00	0.00	A3
ATOM	1154	H	SER	156	33.790	65.577	-2.868	1.00	0.00	A3	ATOM	1205	CA	PIE	161	40.820	63.850	-10.218	1.00	31.21	A3
ATOM	1155	CA	SER	156	34.034	64.105	-4.354	1.00	32.17	A3	ATOM	1206	CB	PIE	161	41.110	65.254	-9.629	1.00	28.28	A3
ATOM	1156	CB	SER	156	32.531	64.319	-4.354	1.00	34.23	A3	ATOM	1207	CG	PIE	161	42.435	65.881	-10.062	1.00	24.92	A3
ATOM	1157	CG	SER	156	32.000	64.195	-5.879	1.00	39.35	A3	ATOM	1208	CD1	PIE	161	43.696	66.228	-11.289	1.00	22.98	A3
ATOM	1158	HC	SER	156	31.120	63.815	-5.851	1.00	0.00	A3	ATOM	1209	CD2	PIE	161	43.464	66.021	-9.135	1.00	23.63	A3
ATOM	1159	C	SER	156	34.845	64.338	-5.632	1.00	33.46	A3	ATOM	1210	CE1	PIE	161	43.941	66.695	-11.767	1.00	21.06	A3
ATOM	1160	O	SER	156	35.411	63.380	-6.174	1.00	34.62	A3	ATOM	1211	CE2	PIE	161	44.701	66.495	-9.528	1.00	20.28	A3
ATOM	1161	N	HIS	157	35.054	65.576	-6.133	1.00	33.90	A3	ATOM	1212	CZ	PIE	161	44.939	66.826	-10.832	1.00	17.01	A3
ATOM	1162	H	HIS	157	34.771	66.349	-5.605	1.00	0.00	A3	ATOM	1213	C	PIE	161	42.008	62.907	-9.943	1.00	31.77	A3
ATOM	1163	CA	HIS	157	35.821	65.773	-7.383	1.00	31.19	A3	ATOM	1214	O	PIE	161	42.786	62.575	-10.845	1.00	32.63	A3
ATOM	1164	CB	HIS	157	35.707	67.209	-7.900	1.00	32.59	A3	ATOM	1215	N	LEU	162	42.117	62.434	-8.690	1.00	31.97	A3
ATOM	1165	CG	HIS	157	34.369	67.449	-8.566	1.00	31.11	A3	ATOM	1216	H	LEU	162	41.420	62.691	-8.054	1.00	0.00	A3
ATOM	1166	CD2	HIS	157	34.127	67.594	-9.928	1.00	30.78	A3	ATOM	1217	CA	LEU	162	43.186	61.574	-8.232	1.00	31.29	A3
ATOM	1167	ND1	HIS	157	33.223	67.666	-7.942	1.00	32.36	A3	ATOM	1218	CB	LEU	162	43.204	61.433	-6.743	1.00	25.84	A3
ATOM	1168	ND1	HIS	157	33.080	67.773	-6.979	1.00	0.00	A3	ATOM	1219	CG	LEU	162	43.604	62.674	-6.003	1.00	26.02	A3
ATOM	1169	CE1	HIS	157	32.293	67.732	-8.875	1.00	32.01	A3	ATOM	1220	CD1	LEU	162	43.594	62.455	-4.516	1.00	25.74	A3
ATOM	1170	NE2	HIS	157	32.838	67.571	-10.060	1.00	29.18	A3	ATOM	1221	CD2	LEU	162	45.107	62.994	-6.415	1.00	27.04	A3
ATOM	1171	HE2	HIS	157	32.327	67.621	-10.895	1.00	0.00	A3	ATOM	1222	C	LEU	162	43.061	60.212	-8.813	1.00	32.21	A3

FIGURE 5

ATOM	1223	O	LEU	162	44.107	59.654	-9.070	1.00	36.51	A3	ATOM	1274	CG1	VAL	168	46.687	60.431	-16.706	1.00	66.78	A3
ATOM	1224	N	GLU	163	41.976	59.589	-9.082	1.00	37.24	A3	ATOM	1275	CG2	VAL	168	48.278	60.879	-14.840	1.00	68.47	A3
ATOM	1225	H	GLU	163	41.072	60.002	-8.826	1.00	0.00	A3	ATOM	1276	C	VAL	168	49.579	58.339	-15.409	1.00	66.45	A3
ATOM	1226	CA	GLU	163	41.975	58.327	-9.771	1.00	40.47	A3	ATOM	1277	O	VAL	168	50.458	58.183	-16.302	1.00	66.22	A3
ATOM	1227	CB	GLU	163	40.566	57.716	-9.835	1.00	45.38	A3	ATOM	1278	N	LEU	169	49.823	58.241	-14.177	1.00	68.83	A3
ATOM	1228	CG	GLU	163	40.264	56.975	-8.526	1.00	51.84	A3	ATOM	1279	H	LEU	169	49.102	58.404	-13.536	1.00	0.00	A3
ATOM	1229	CD	GLU	163	41.291	55.889	-8.126	1.00	57.97	A3	ATOM	1280	CA	LEU	169	51.141	57.899	-13.695	1.00	71.81	A3
ATOM	1230	OE1	GLU	163	40.897	54.722	-8.092	1.00	61.01	A3	ATOM	1281	CB	LEU	169	51.249	58.228	-12.188	1.00	71.51	A3
ATOM	1231	OE2	GLU	163	42.466	56.180	-7.832	1.00	59.17	A3	ATOM	1282	CG	LEU	169	51.137	59.732	-11.813	1.00	70.68	A3
ATOM	1232	C	GLU	163	42.586	58.430	-11.142	1.00	41.34	A3	ATOM	1283	CD1	LEU	169	51.187	59.876	-10.298	1.00	69.39	A3
ATOM	1233	O	GLU	163	43.456	57.633	-11.486	1.00	42.17	A3	ATOM	1284	CD2	LEU	169	52.223	60.580	-12.491	1.00	68.49	A3
ATOM	1234	N	VAL	164	42.257	59.436	-11.920	1.00	42.28	A3	ATOM	1285	C	LEU	169	51.333	56.414	-13.979	1.00	73.61	A3
ATOM	1235	H	VAL	164	41.589	60.091	-11.615	1.00	0.00	A3	ATOM	1286	O	LEU	169	52.408	56.013	-14.429	1.00	74.75	A3
ATOM	1236	CA	VAL	164	42.911	59.609	-13.187	1.00	44.13	A3	ATOM	1287	N	ARG	170	50.309	55.583	-13.819	1.00	75.45	A3
ATOM	1237	CB	VAL	164	42.707	60.711	-13.940	1.00	45.52	A3	ATOM	1288	H	ARG	170	49.488	55.923	-13.399	1.00	0.00	A3
ATOM	1238	CG1	VAL	164	42.892	60.975	-15.278	1.00	48.79	A3	ATOM	1289	CA	ARG	170	50.364	54.179	-14.199	1.00	78.17	A3
ATOM	1239	CG2	VAL	164	40.786	60.269	-14.226	1.00	46.09	A3	ATOM	1290	CB	ARG	170	48.944	53.642	-14.004	1.00	78.45	A3
ATOM	1240	C	VAL	164	44.386	59.933	-12.991	1.00	46.13	A3	ATOM	1291	CG	ARG	170	48.394	52.506	-14.871	1.00	78.17	A3
ATOM	1241	O	VAL	164	45.192	59.473	-13.794	1.00	45.99	A3	ATOM	1292	CD	ARG	170	48.744	51.181	-14.271	1.00	77.25	A3
ATOM	1242	N	SER	165	44.879	60.677	-12.006	1.00	49.51	A3	ATOM	1293	HE	ARG	170	48.123	51.120	-12.970	1.00	76.15	A3
ATOM	1243	H	SER	165	44.287	61.173	-11.396	1.00	0.00	A3	ATOM	1294	HE	ARG	170	47.245	51.528	-12.824	1.00	0.00	A3
ATOM	1244	CA	SER	165	46.325	60.845	-11.895	1.00	53.44	A3	ATOM	1295	CZ	ARG	170	48.758	50.547	-11.970	1.00	76.14	A3
ATOM	1245	CB	SER	165	46.715	61.796	-10.775	1.00	54.77	A3	ATOM	1296	NH1	ARG	170	49.973	50.017	-12.112	1.00	76.84	A3
ATOM	1246	CG	SER	165	46.049	61.618	-9.530	1.00	59.99	A3	ATOM	1297	NH11	ARG	170	50.441	50.030	-12.994	1.00	0.00	A3
ATOM	1247	HG	SER	165	45.997	60.694	-9.261	1.00	0.00	A3	ATOM	1298	NH12	ARG	170	50.406	49.570	-11.329	1.00	0.00	A3
ATOM	1248	C	SER	165	46.958	59.502	-11.630	1.00	55.15	A3	ATOM	1299	NH2	ARG	170	48.147	50.492	-10.806	1.00	77.02	A3
ATOM	1249	O	SER	165	48.028	59.227	-12.148	1.00	55.02	A3	ATOM	1300	NH21	ARG	170	47.237	50.890	-10.714	1.00	0.00	A3
ATOM	1250	N	TYR	166	46.239	58.645	-10.900	1.00	58.57	A3	ATOM	1301	NH22	ARG	170	48.586	50.052	-10.023	1.00	0.00	A3
ATOM	1251	H	TYR	166	45.324	58.948	-10.549	1.00	0.00	A3	ATOM	1302	C	ARG	170	50.870	54.052	-15.647	1.00	79.84	A3
ATOM	1252	CA	TYR	166	46.617	57.273	-10.625	1.00	61.42	A3	ATOM	1303	O	ARG	170	51.924	53.470	-15.908	1.00	80.07	A3
ATOM	1253	CB	TYR	166	45.543	56.653	-9.680	1.00	64.05	A3	ATOM	1304	N	HIS	171	50.193	54.663	-16.611	1.00	81.58	A3
ATOM	1254	CG	TYR	166	45.502	55.138	-9.682	1.00	69.00	A3	ATOM	1305	H	HIS	171	49.433	55.234	-16.359	1.00	0.00	A3
ATOM	1255	CD1	TYR	166	44.389	54.501	-10.185	1.00	71.64	A3	ATOM	1306	CA	HIS	171	50.663	54.597	-17.970	1.00	84.03	A3
ATOM	1256	CE1	TYR	166	44.367	53.130	-10.283	1.00	73.15	A3	ATOM	1307	CB	HIS	171	49.590	55.054	-18.902	1.00	86.82	A3
ATOM	1257	CD2	TYR	166	46.594	54.409	-9.257	1.00	71.27	A3	ATOM	1308	CG	HIS	171	48.496	54.037	-19.147	1.00	90.73	A3
ATOM	1258	CE2	TYR	166	46.584	53.040	-9.346	1.00	72.92	A3	ATOM	1309	CD2	HIS	171	47.467	53.765	-18.272	1.00	91.35	A3
ATOM	1259	CZ	TYR	166	45.468	52.417	-9.862	1.00	75.71	A3	ATOM	1310	ND1	HIS	171	48.308	53.301	-20.248	1.00	92.24	A3
ATOM	1260	OH	TYR	166	45.474	51.038	-10.016	1.00	80.61	A3	ATOM	1311	ND1	HIS	171	48.887	53.287	-21.044	1.00	0.00	A3
ATOM	1261	H11	TYR	166	46.571	50.736	-10.134	1.00	0.00	A3	ATOM	1312	CE1	HIS	171	47.204	52.605	-20.077	1.00	92.41	A3
ATOM	1262	C	TYR	166	46.712	56.567	-11.987	1.00	62.34	A3	ATOM	1313	NE2	HIS	171	46.711	52.892	-18.891	1.00	92.59	A3
ATOM	1263	O	TYR	166	47.766	55.981	-12.282	1.00	63.25	A3	ATOM	1314	NE2	HIS	171	45.884	52.511	-18.518	1.00	0.00	A3
ATOM	1264	N	ALA	167	45.727	56.622	-12.884	1.00	61.27	A3	ATOM	1315	C	HIS	171	51.907	55.446	-18.232	1.00	85.42	A3
ATOM	1265	H	ALA	167	44.893	57.089	-12.678	1.00	0.00	A3	ATOM	1316	O	HIS	171	52.440	55.352	-19.344	1.00	85.98	A3
ATOM	1266	CA	ALA	167	45.933	55.982	-14.159	1.00	61.47	A3	ATOM	1317	N	LEU	172	52.359	56.307	-17.302	1.00	86.13	A3
ATOM	1267	CB	ALA	167	46.608	55.904	-14.904	1.00	60.98	A3	ATOM	1318	H	LEU	172	51.870	56.411	-16.463	1.00	0.00	A3
ATOM	1268	C	ALA	167	46.982	56.694	-15.070	1.00	62.19	A3	ATOM	1319	CA	LEU	172	53.550	57.133	-17.496	1.00	86.02	A3
ATOM	1269	O	ALA	167	47.719	56.000	-15.734	1.00	62.63	A3	ATOM	1320	CB	LEU	172	53.500	58.357	-16.607	1.00	86.31	A3
ATOM	1270	N	VAL	168	47.210	58.011	-14.991	1.00	63.37	A3	ATOM	1321	CG	LEU	172	54.022	59.658	-17.203	1.00	87.48	A3
ATOM	1271	H	VAL	168	46.756	58.570	-14.330	1.00	0.00	A3	ATOM	1322	CD1	LEU	172	53.436	59.939	-18.596	1.00	87.68	A3
ATOM	1272	CA	VAL	168	48.174	58.593	-15.923	1.00	65.62	A3	ATOM	1323	CD2	LEU	172	53.645	60.778	-16.251	1.00	87.95	A3
ATOM	1273	CB	VAL	168	48.061	60.121	-16.131	1.00	66.30	A3	ATOM	1324	C	LEU	172	54.813	56.357	-17.180	1.00	85.92	A3

FIGURE 5

ATOM 1325 O LEU 172	55.896	56.660	-17.692	1.00	86.23	A3	ATOM 1376 CD1 PHE 214	45.176	39.459	23.044	1.00	-42.77	BI
ATOM 1326 N ALA 173	54.733	55.383	-16.282	1.00	85.49	A3	ATOM 1377 CD2 PHE 214	46.818	37.794	23.400	1.00	-42.02	BI
ATOM 1327 H ALA 173	53.899	55.276	-15.769	1.00	0.00	A3	ATOM 1378 CE1 PHE 214	44.197	38.554	23.423	1.00	-41.82	BI
ATOM 1328 CA ALA 173	55.856	54.497	-16.087	1.00	85.65	A3	ATOM 1379 CE2 PHE 214	45.834	36.898	23.776	1.00	-41.70	BI
ATOM 1329 CB ALA 173	56.602	54.859	-14.809	1.00	85.01	A3	ATOM 1380 CZ PHE 214	44.519	37.277	23.791	1.00	-41.05	BI
ATOM 1330 C ALA 173	55.330	53.073	-16.008	1.00	86.54	A3	ATOM 1381 C PHE 214	47.109	39.656	20.321	1.00	16.54	BI
ATOM 1331 OT1 ALA 173	55.585	52.347	-16.971	1.00	87.21	A3	ATOM 1382 O PHE 214	46.735	38.566	19.889	1.00	37.99	BI
ATOM 1332 OT2 ALA 173	54.650	52.707	-15.036	1.00	87.31	A3	ATOM 1383 N LEU 215	46.616	40.812	19.893	1.00	33.27	BI
ATOM 1333 CB LEU 210	45.234	42.591	25.453	1.00	52.47	BI	ATOM 1384 H LEU 215	47.008	41.642	20.238	1.00	0.00	BI
ATOM 1334 CG LEU 210	43.799	42.058	25.547	1.00	51.68	BI	ATOM 1385 CA LEU 215	45.504	40.864	18.966	1.00	30.38	BI
ATOM 1335 CD1 LEU 210	43.123	42.562	26.804	1.00	53.37	BI	ATOM 1386 CB LEU 215	45.099	42.282	18.701	1.00	31.82	BI
ATOM 1336 CD2 LEU 210	43.050	42.453	24.303	1.00	51.37	BI	ATOM 1387 CG LEU 215	43.857	42.530	17.893	1.00	32.78	BI
ATOM 1337 C LEU 210	46.770	44.374	24.596	1.00	50.98	BI	ATOM 1388 CD1 LEU 215	42.727	41.963	18.737	1.00	32.95	BI
ATOM 1338 O LEU 210	46.475	45.767	23.790	1.00	51.76	BI	ATOM 1389 CD2 LEU 215	43.688	44.011	17.508	1.00	28.93	BI
ATOM 1339 HT1 LEU 210	44.382	44.922	24.421	1.00	0.00	BI	ATOM 1390 C LEU 215	45.811	40.232	17.648	1.00	29.57	BI
ATOM 1340 HT2 LEU 210	45.157	45.974	25.414	1.00	0.00	BI	ATOM 1391 O LEU 215	44.922	39.632	17.055	1.00	31.28	BI
ATOM 1341 N LEU 210	44.705	45.041	25.406	1.00	53.59	BI	ATOM 1392 N LEU 216	47.031	40.379	17.155	1.00	29.44	BI
ATOM 1342 HT3 LEU 210	43.855	45.012	25.997	1.00	0.00	BI	ATOM 1393 H LEU 216	47.677	40.935	17.646	1.00	0.00	BI
ATOM 1343 CA LEU 210	45.730	44.038	25.676	1.00	52.35	BI	ATOM 1394 CA LEU 216	47.465	39.790	15.893	1.00	29.89	BI
ATOM 1344 N PRO 211	47.974	43.825	24.494	1.00	49.35	BI	ATOM 1395 CB LEU 216	48.791	40.450	15.472	1.00	28.61	BI
ATOM 1345 CD PRO 211	48.621	43.024	25.532	1.00	49.52	BI	ATOM 1396 CG LEU 216	48.682	41.877	14.919	1.00	26.84	BI
ATOM 1346 CA PRO 211	48.895	44.191	23.419	1.00	49.04	BI	ATOM 1397 CD1 LEU 216	49.925	42.558	15.344	1.00	28.57	BI
ATOM 1347 CB PRO 211	50.209	43.571	23.865	1.00	49.02	BI	ATOM 1398 CD2 LEU 216	48.446	41.950	13.452	1.00	24.09	BI
ATOM 1348 CG PRO 211	49.794	42.438	24.783	1.00	49.77	BI	ATOM 1399 C LEU 216	47.613	38.274	16.062	1.00	31.23	BI
ATOM 1349 C PRO 211	48.543	43.864	21.965	1.00	48.03	BI	ATOM 1400 O LEU 216	47.328	37.514	15.138	1.00	29.20	BI
ATOM 1350 O PRO 211	47.872	42.896	21.622	1.00	49.05	BI	ATOM 1401 N LYS 217	47.999	37.826	17.261	1.00	32.50	BI
ATOM 1351 N GLN 212	49.032	44.675	21.051	1.00	46.52	BI	ATOM 1402 H LYS 217	48.305	38.482	17.926	1.00	0.00	BI
ATOM 1352 H GLN 212	49.506	45.478	21.349	1.00	0.00	BI	ATOM 1403 CA LYS 217	48.067	36.439	17.599	1.00	34.90	BI
ATOM 1353 CA GLN 212	48.839	44.461	19.641	1.00	45.47	BI	ATOM 1404 CB LYS 217	48.645	36.280	19.002	1.00	38.07	BI
ATOM 1354 CB GLN 212	49.533	45.522	18.849	1.00	46.81	BI	ATOM 1405 CG LYS 217	49.394	34.978	19.109	1.00	45.25	BI
ATOM 1355 CG GLN 212	48.482	46.139	17.999	1.00	49.55	BI	ATOM 1406 CD LYS 217	49.714	34.491	20.521	1.00	53.27	BI
ATOM 1356 CD GLN 212	49.024	46.703	16.709	1.00	54.21	BI	ATOM 1407 CE LYS 217	50.229	33.024	20.297	1.00	59.03	BI
ATOM 1357 OE1 GLN 212	48.429	47.672	16.232	1.00	57.72	BI	ATOM 1408 NZ LYS 217	50.213	32.135	21.467	1.00	62.10	BI
ATOM 1358 NE2 GLN 212	50.086	46.176	16.074	1.00	52.39	BI	ATOM 1409 H21 LYS 217	49.239	32.056	21.824	1.00	0.00	BI
ATOM 1359 HE1 GLN 212	50.530	45.383	16.430	1.00	0.00	BI	ATOM 1410 H22 LYS 217	50.830	32.515	22.214	1.00	0.00	BI
ATOM 1360 HE2 GLN 212	50.341	46.625	15.244	1.00	0.00	BI	ATOM 1411 H23 LYS 217	50.554	31.195	21.179	1.00	0.00	BI
ATOM 1361 C GLN 212	49.390	43.139	19.185	1.00	44.79	BI	ATOM 1412 C LYS 217	46.617	35.950	17.546	1.00	36.77	BI
ATOM 1362 O GLN 212	48.959	42.520	18.208	1.00	44.01	BI	ATOM 1413 O LYS 217	46.311	34.933	16.886	1.00	39.58	BI
ATOM 1363 N SER 213	50.401	42.671	19.893	1.00	44.72	BI	ATOM 1414 N CYS 218	45.664	36.638	18.177	1.00	34.86	BI
ATOM 1364 H SER 213	50.730	43.115	20.698	1.00	0.00	BI	ATOM 1415 H CYS 218	45.907	37.388	18.751	1.00	0.00	BI
ATOM 1365 CA SER 213	51.025	41.424	19.521	1.00	43.76	BI	ATOM 1416 CA CYS 218	44.277	36.238	18.076	1.00	33.61	BI
ATOM 1366 CB SER 213	52.220	41.124	20.354	1.00	45.29	BI	ATOM 1417 CB CYS 218	43.430	37.175	18.846	1.00	33.71	BI
ATOM 1367 CG SER 213	51.802	41.455	21.681	1.00	52.50	BI	ATOM 1418 SG CYS 218	43.856	36.710	20.515	1.00	35.92	BI
ATOM 1368 HG SER 213	52.479	41.127	22.288	1.00	0.00	BI	ATOM 1419 C CYS 218	43.766	36.189	16.632	1.00	32.89	BI
ATOM 1369 C SER 213	50.014	40.376	19.784	1.00	40.92	BI	ATOM 1420 O CYS 218	43.155	35.169	16.323	1.00	34.71	BI
ATOM 1370 O SER 213	49.964	39.492	18.947	1.00	43.32	BI	ATOM 1421 H LEU 219	44.035	37.169	15.777	1.00	29.52	BI
ATOM 1371 N PHE 214	49.242	40.571	20.876	1.00	38.86	BI	ATOM 1422 H LEU 219	43.512	37.960	16.104	1.00	0.00	BI
ATOM 1372 H PHE 214	49.414	41.370	21.410	1.00	0.00	BI	ATOM 1423 CA LEU 219	43.614	37.119	14.393	1.00	27.44	BI
ATOM 1373 CA PHE 214	48.210	39.664	21.336	1.00	37.40	BI	ATOM 1424 CB LEU 219	44.116	38.412	13.727	1.00	26.24	BI
ATOM 1374 CB PHE 214	47.568	40.064	22.634	1.00	17.45	BI	ATOM 1425 CG LEU 219	43.884	38.768	12.241	1.00	25.07	BI
ATOM 1375 CG PHE 214	46.494	39.080	23.035	1.00	41.01	BI	ATOM 1426 CD LEU 219	42.402	38.975	11.996	1.00	24.24	BI

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ATOM	1427	CD2	LEU	219	44.563	40.051	11.882	1.00	22.10	BI	ATOM	1478	II	LYS	224	42.791	31.235	12.048	1.00	0.00	BI
ATOM	1428	C	LEU	219	44.121	35.867	13.634	1.00	28.14	BI	ATOM	1479	CA	LYS	224	42.714	29.411	10.994	1.00	27.70	BI
ATOM	1429	O	LEU	219	43.373	35.204	12.889	1.00	27.12	BI	ATOM	1480	CB	LYS	224	43.922	29.085	11.818	1.00	30.07	BI
ATOM	1430	N	GLU	220	45.399	35.499	13.795	1.00	28.06	BI	ATOM	1481	CG	LYS	224	44.372	27.660	11.706	1.00	36.70	BI
ATOM	1431	II	GLU	220	45.957	35.974	14.448	1.00	0.00	BI	ATOM	1482	CD	LYS	224	45.829	27.544	12.127	1.00	41.68	BI
ATOM	1432	CA	GLU	220	45.963	34.411	13.048	1.00	28.38	BI	ATOM	1483	CE	LYS	224	46.303	26.478	11.131	1.00	48.18	BI
ATOM	1433	CB	GLU	220	47.376	34.198	13.469	1.00	34.25	BI	ATOM	1484	NZ	LYS	224	47.750	26.492	10.913	1.00	53.57	BI
ATOM	1434	CG	GLU	220	48.049	33.079	12.666	1.00	46.36	BI	ATOM	1485	IZ1	LYS	224	48.230	26.241	11.801	1.00	0.00	BI
ATOM	1435	CD	GLU	220	49.345	32.794	12.907	1.00	55.51	BI	ATOM	1486	IZ2	LYS	224	48.057	27.436	10.606	1.00	0.00	BI
ATOM	1436	OE1	GLU	220	50.113	32.133	12.021	1.00	58.95	BI	ATOM	1487	IZ3	LYS	224	47.998	25.792	10.183	1.00	0.00	BI
ATOM	1437	OE2	GLU	220	50.144	33.213	13.930	1.00	60.41	BI	ATOM	1488	C	LYS	224	41.464	28.598	11.347	1.00	26.27	BI
ATOM	1438	C	GLU	220	45.134	33.193	13.354	1.00	27.30	BI	ATOM	1489	O	LYS	224	40.970	27.810	10.510	1.00	24.82	BI
ATOM	1439	O	GLU	220	44.662	32.524	12.437	1.00	27.08	BI	ATOM	1490	N	ILE	225	40.892	28.835	12.547	1.00	24.75	BI
ATOM	1440	N	GLN	221	44.866	33.023	14.642	1.00	25.42	BI	ATOM	1491	II	ILE	225	41.308	29.487	13.151	1.00	0.00	BI
ATOM	1441	II	GLN	221	45.229	33.687	15.268	1.00	0.00	BI	ATOM	1492	CA	ILE	225	39.636	28.147	12.943	1.00	23.33	BI
ATOM	1442	CA	GLN	221	44.074	31.940	15.176	1.00	26.28	BI	ATOM	1493	CB	ILE	225	39.146	28.622	14.296	1.00	18.08	BI
ATOM	1443	CB	GLN	221	44.143	31.927	16.691	1.00	26.78	BI	ATOM	1494	CG2	ILE	225	37.874	27.872	14.577	1.00	15.43	BI
ATOM	1444	CG	GLN	221	45.555	31.456	17.011	1.00	29.19	BI	ATOM	1495	CD	ILE	225	40.161	28.400	15.380	1.00	13.18	BI
ATOM	1445	CD	GLN	221	45.752	31.067	18.442	1.00	31.98	BI	ATOM	1496	CE	ILE	225	39.787	28.967	16.749	1.00	13.57	BI
ATOM	1446	OE1	GLN	221	46.472	30.162	18.808	1.00	35.98	BI	ATOM	1497	C	ILE	225	38.594	28.437	11.889	1.00	27.28	BI
ATOM	1447	HE21	GLN	221	45.110	31.736	19.347	1.00	39.31	BI	ATOM	1498	O	ILE	225	37.978	27.492	11.400	1.00	31.49	BI
ATOM	1448	HE21	GLN	221	45.263	31.423	20.246	1.00	0.00	BI	ATOM	1499	N	GLN	226	38.396	29.677	11.402	1.00	29.69	BI
ATOM	1449	HE22	GLN	221	44.571	32.514	19.111	1.00	0.00	BI	ATOM	1500	II	GLN	226	38.894	30.413	11.803	1.00	0.00	BI
ATOM	1450	C	GLN	221	42.615	31.925	14.789	1.00	26.21	BI	ATOM	1501	CA	GLN	226	37.450	29.969	10.313	1.00	29.12	BI
ATOM	1451	O	GLN	221	42.186	30.896	14.269	1.00	30.69	BI	ATOM	1502	CB	GLN	226	37.366	31.438	9.962	1.00	32.26	BI
ATOM	1452	N	VAL	222	41.814	32.962	14.984	1.00	23.63	BI	ATOM	1503	CG	GLN	226	36.682	32.156	11.108	1.00	46.28	BI
ATOM	1453	II	VAL	222	42.199	33.746	15.426	1.00	0.00	BI	ATOM	1504	CD	GLN	226	36.429	33.613	10.816	1.00	37.88	BI
ATOM	1454	CA	VAL	222	40.429	33.034	14.537	1.00	21.92	BI	ATOM	1505	OE1	GLN	226	37.158	34.281	10.076	1.00	36.34	BI
ATOM	1455	CB	VAL	222	39.934	34.442	14.793	1.00	21.36	BI	ATOM	1506	NE2	GLN	226	35.359	34.114	11.421	1.00	39.62	BI
ATOM	1456	CG1	VAL	222	38.706	34.831	14.027	1.00	17.72	BI	ATOM	1507	HE21	GLN	226	34.823	33.501	11.971	1.00	0.00	BI
ATOM	1457	CG2	VAL	222	39.671	34.496	16.257	1.00	20.95	BI	ATOM	1508	HE22	GLN	226	35.153	35.057	11.287	1.00	0.00	BI
ATOM	1458	C	VAL	222	40.374	32.707	13.066	1.00	22.65	BI	ATOM	1509	C	GLN	226	37.714	29.295	9.007	1.00	26.82	BI
ATOM	1459	O	VAL	222	39.475	32.013	12.632	1.00	23.72	BI	ATOM	1510	O	GLN	226	36.775	28.887	8.325	1.00	27.45	BI
ATOM	1460	N	ARG	223	41.341	33.120	12.283	1.00	23.95	BI	ATOM	1511	N	GLY	227	38.940	29.186	8.570	1.00	26.55	BI
ATOM	1461	II	ARG	223	42.099	33.614	12.666	1.00	0.00	BI	ATOM	1512	II	GLY	227	39.688	29.612	9.043	1.00	0.00	BI
ATOM	1462	CA	ARG	223	41.309	32.939	10.844	1.00	27.19	BI	ATOM	1513	CA	GLY	227	39.195	28.427	7.348	1.00	27.27	BI
ATOM	1463	CB	ARG	223	42.794	33.935	10.783	1.00	29.26	BI	ATOM	1514	C	GLY	227	38.832	26.949	7.574	1.00	27.65	BI
ATOM	1464	CG	ARG	223	42.102	34.364	8.869	1.00	35.23	BI	ATOM	1515	O	GLY	227	38.287	26.291	6.656	1.00	26.79	BI
ATOM	1465	CD	ARG	223	42.880	33.487	7.929	1.00	41.88	BI	ATOM	1516	N	ASP	228	39.025	26.429	8.819	1.00	27.03	BI
ATOM	1466	NE	ARG	223	41.972	32.676	7.096	1.00	47.42	BI	ATOM	1517	II	ASP	228	39.460	26.957	9.523	1.00	0.00	BI
ATOM	1467	IE	ARG	223	41.451	31.953	7.502	1.00	0.00	BI	ATOM	1518	CA	ASP	228	38.618	25.038	9.052	1.00	28.20	BI
ATOM	1468	CZ	ARG	223	41.875	32.896	5.784	1.00	46.15	BI	ATOM	1519	CB	ASP	228	38.986	24.492	10.391	1.00	26.04	BI
ATOM	1469	NI1	ARG	223	42.575	33.837	5.246	1.00	49.54	BI	ATOM	1520	CG	ASP	228	40.427	24.554	10.774	1.00	24.88	BI
ATOM	1470	HI11	ARG	223	42.522	33.989	4.259	1.00	0.00	BI	ATOM	1521	OD1	ASP	228	40.627	24.521	11.977	1.00	23.37	BI
ATOM	1471	HI12	ARG	223	43.156	34.428	5.805	1.00	0.00	BI	ATOM	1522	OD2	ASP	228	41.302	24.637	9.912	1.00	23.32	BI
ATOM	1472	NI12	ARG	223	41.178	32.161	4.952	1.00	45.74	BI	ATOM	1523	C	ASP	228	37.120	24.830	8.992	1.00	27.23	BI
ATOM	1473	HI121	ARG	223	40.697	31.353	5.290	1.00	0.00	BI	ATOM	1524	O	ASP	228	36.662	23.900	8.336	1.00	27.07	BI
ATOM	1474	HI122	ARG	223	41.154	32.399	3.980	1.00	0.00	BI	ATOM	1525	N	GLY	229	36.390	25.739	9.639	1.00	26.74	BI
ATOM	1475	C	ARG	223	41.624	31.492	10.430	1.00	29.13	BI	ATOM	1526	II	GLY	229	36.861	26.444	10.134	1.00	0.00	BI
ATOM	1476	O	ARG	223	41.181	30.987	9.376	1.00	29.32	BI	ATOM	1527	CA	GLY	229	34.946	25.723	9.673	1.00	25.87	BI
ATOM	1477	N	LYS	224	42.413	30.791	11.259	1.00	29.17	BI	ATOM	1528	C	GLY	229	34.593	25.825	8.274	1.00	24.95	BI

FIGURE 5

ATOM	1631	N	LVS	241	24.174	18.011	0.694	1.00	37.36	BI	ATOM	1682	CA	GIU	246	19.257	29.229	7.731	1.00	41.10	BI
ATOM	1632	II	LVS	241	25.091	18.023	0.345	1.00	0.00	BI	ATOM	1683	CB	GIU	246	19.044	30.107	6.438	1.00	41.15	BI
ATOM	1633	CA	LVS	241	23.314	19.115	0.275	1.00	36.37	BI	ATOM	1684	CG	GIU	246	20.256	30.918	5.944	1.00	47.07	BI
ATOM	1634	CB	LVS	241	22.173	18.648	-0.595	1.00	38.38	BI	ATOM	1685	CD	GIU	246	20.813	30.539	4.358	1.00	52.63	BI
ATOM	1635	CG	LVS	241	22.645	17.940	-1.838	1.00	42.94	BI	ATOM	1686	OEI	GIU	246	22.054	30.545	4.374	1.00	54.22	BI
ATOM	1636	CD	LVS	241	23.468	18.809	-2.737	1.00	46.97	BI	ATOM	1687	OEZ	GIU	246	20.002	30.250	3.656	1.00	54.39	BI
ATOM	1637	CE	LVS	241	23.657	18.070	-4.051	1.00	49.20	BI	ATOM	1688	C	GIU	246	18.071	28.298	7.819	1.00	40.57	BI
ATOM	1638	NZ	LVS	241	22.509	18.372	-4.893	1.00	51.34	BI	ATOM	1689	O	GIU	246	17.308	28.338	8.791	1.00	39.90	BI
ATOM	1639	IZ1	LVS	241	22.447	19.400	-5.038	1.00	0.00	BI	ATOM	1690	N	GIU	247	18.025	27.388	6.840	1.00	40.32	BI
ATOM	1640	IZ2	LVS	241	21.641	18.041	-4.426	1.00	0.00	BI	ATOM	1691	II	GIU	247	18.750	27.334	6.190	1.00	0.00	BI
ATOM	1641	IZ3	LVS	241	22.609	17.895	-5.811	1.00	0.00	BI	ATOM	1692	CA	GIU	247	17.001	26.347	6.830	1.00	40.76	BI
ATOM	1642	C	LVS	241	22.720	19.904	1.429	1.00	33.37	BI	ATOM	1693	CB	GIU	247	17.139	25.423	5.642	1.00	44.03	BI
ATOM	1643	O	LVS	241	21.728	20.580	1.223	1.00	31.90	BI	ATOM	1694	CG	GIU	247	16.830	26.240	4.400	1.00	48.34	BI
ATOM	1644	N	LEU	242	23.286	19.853	2.648	1.00	31.40	BI	ATOM	1695	CD	GIU	247	17.163	25.628	3.050	1.00	50.24	BI
ATOM	1645	H	LEU	242	24.055	19.260	2.756	1.00	0.00	BI	ATOM	1696	OEI	GIU	247	16.849	26.299	2.056	1.00	52.92	BI
ATOM	1646	CA	LEU	242	22.904	20.682	3.758	1.00	31.09	BI	ATOM	1697	OEZ	GIU	247	17.744	24.533	2.987	1.00	50.84	BI
ATOM	1647	CB	LEU	242	23.253	20.059	5.096	1.00	28.55	BI	ATOM	1698	C	GIU	247	16.966	25.444	8.034	1.00	59.24	BI
ATOM	1648	CG	LEU	242	22.571	18.798	5.641	1.00	30.36	BI	ATOM	1699	O	GIU	247	15.915	24.888	8.329	1.00	39.40	BI
ATOM	1649	CD1	LEU	242	22.530	18.814	7.138	1.00	29.62	BI	ATOM	1700	N	LEU	248	18.066	25.280	8.760	1.00	37.92	BI
ATOM	1650	CD2	LEU	242	21.086	18.861	3.543	1.00	31.94	BI	ATOM	1701	II	LEU	248	18.864	25.814	8.576	1.00	0.00	BI
ATOM	1651	C	LEU	242	23.778	21.933	3.550	1.00	34.03	BI	ATOM	1702	CA	LEU	248	18.101	24.338	9.858	1.00	35.75	BI
ATOM	1652	O	LEU	242	24.903	22.027	4.058	1.00	35.53	BI	ATOM	1703	CB	LEU	248	19.458	23.623	9.796	1.00	34.11	BI
ATOM	1653	N	CYS	243	23.316	22.883	2.722	1.00	34.89	BI	ATOM	1704	CG	LEU	248	19.669	22.866	8.430	1.00	34.00	BI
ATOM	1654	H	CYS	243	22.491	22.663	2.238	1.00	0.00	BI	ATOM	1705	CD1	LEU	248	20.997	22.149	8.306	1.00	33.97	BI
ATOM	1655	CA	CYS	243	24.051	24.083	2.377	1.00	35.42	BI	ATOM	1706	CD2	LEU	248	18.620	21.810	8.322	1.00	32.33	BI
ATOM	1656	C	CYS	243	23.492	25.335	2.975	1.00	36.85	BI	ATOM	1707	C	LEU	248	17.871	25.031	11.153	1.00	36.51	BI
ATOM	1657	O	CYS	243	23.956	26.400	2.565	1.00	40.10	BI	ATOM	1708	N	LEU	248	17.736	24.370	12.186	1.00	36.31	BI
ATOM	1658	CB	CYS	243	24.046	24.383	0.979	1.00	33.12	BI	ATOM	1709	II	VAL	249	17.663	26.350	11.146	1.00	38.88	BI
ATOM	1659	SG	CYS	243	24.438	22.883	0.099	1.00	38.25	BI	ATOM	1710	II	VAL	249	17.566	26.810	10.281	1.00	0.00	BI
ATOM	1660	N	IIS	244	22.496	25.393	3.848	1.00	35.37	BI	ATOM	1711	CA	VAL	249	17.573	27.133	12.371	1.00	41.39	BI
ATOM	1661	H	IIS	244	22.185	24.588	4.318	1.00	0.00	BI	ATOM	1712	CB	VAL	249	17.265	28.640	12.020	1.00	43.72	BI
ATOM	1662	CA	IIS	244	21.939	26.676	4.191	1.00	33.29	BI	ATOM	1713	CG1	VAL	249	15.804	28.985	11.776	1.00	44.70	BI
ATOM	1663	CB	IIS	244	20.655	26.987	3.340	1.00	33.64	BI	ATOM	1714	CG2	VAL	249	17.702	29.434	13.214	1.00	45.20	BI
ATOM	1664	CG	IIS	244	20.915	27.205	1.857	1.00	33.12	BI	ATOM	1715	C	VAL	249	16.590	26.635	13.406	1.00	42.01	BI
ATOM	1665	CD2	IIS	244	20.288	26.584	0.814	1.00	37.29	BI	ATOM	1716	O	VAL	249	16.912	26.716	14.594	1.00	44.77	BI
ATOM	1666	ND1	IIS	244	21.874	27.902	1.298	1.00	36.85	BI	ATOM	1717	N	LEU	250	15.453	26.035	13.016	1.00	41.61	BI
ATOM	1667	IDI	IIS	244	22.648	28.281	1.778	1.00	0.00	BI	ATOM	1718	II	LEU	250	15.319	25.919	12.053	1.00	0.00	BI
ATOM	1668	CE1	IIS	244	21.874	27.722	-0.013	1.00	35.95	BI	ATOM	1719	CA	LEU	250	14.457	25.537	13.987	1.00	43.96	BI
ATOM	1669	NE2	IIS	244	20.910	26.920	-0.301	1.00	35.54	BI	ATOM	1720	CB	LEU	250	13.102	25.296	13.373	1.00	41.88	BI
ATOM	1670	HE2	IIS	244	20.616	26.706	-1.214	1.00	0.00	BI	ATOM	1721	CG	LEU	250	12.729	26.781	12.313	1.00	47.04	BI
ATOM	1671	C	IIS	244	21.621	26.565	5.650	1.00	33.38	BI	ATOM	1722	CD1	LEU	250	13.092	25.577	11.011	1.00	47.60	BI
ATOM	1672	O	IIS	244	20.546	26.105	6.029	1.00	33.23	BI	ATOM	1723	CD2	LEU	250	11.286	26.772	12.441	1.00	46.18	BI
ATOM	1673	N	PRO	245	22.539	27.018	6.499	1.00	33.21	BI	ATOM	1724	C	LEU	250	14.832	24.207	14.626	1.00	43.96	BI
ATOM	1674	CD	PRO	245	23.851	27.524	6.099	1.00	31.29	BI	ATOM	1725	O	LEU	250	14.450	23.887	15.764	1.00	41.07	BI
ATOM	1675	CA	PRO	245	22.373	26.979	7.948	1.00	34.16	BI	ATOM	1726	N	LEU	251	15.691	23.446	13.893	1.00	42.41	BI
ATOM	1676	CB	PRO	245	23.490	27.799	8.467	1.00	32.85	BI	ATOM	1727	II	LEU	251	16.049	23.788	13.048	1.00	0.00	BI
ATOM	1677	CG	PRO	245	24.564	27.549	7.428	1.00	31.74	BI	ATOM	1728	CA	LEU	251	16.155	22.159	14.362	1.00	40.63	BI
ATOM	1678	C	PRO	245	21.032	27.470	8.407	1.00	36.26	BI	ATOM	1729	CB	LEU	251	16.834	21.418	13.257	1.00	36.77	BI
ATOM	1679	O	PRO	245	20.478	26.878	9.315	1.00	38.13	BI	ATOM	1730	CG	LEU	251	15.996	20.629	12.267	1.00	33.76	BI
ATOM	1680	N	GIU	246	20.529	28.463	7.640	1.00	39.64	BI	ATOM	1731	CD1	LEU	251	14.595	21.168	11.956	1.00	33.46	BI
ATOM	1681	II	GIU	246	21.134	28.747	6.934	1.00	0.00	BI	ATOM	1732	CD2	LEU	251	16.875	20.619	11.050	1.00	44.71	BI

FIGURE 5

ATOM 1529 O GLY 229	33.370	25.222	7.956	1.00	25.73	BI	ATOM 1580 NZ LYS 235	37.117	17.460	2.521	1.00	-41.44	BI
ATOM 1530 N ALA 230	35.058	26.541	7.391	1.00	23.97	BI	ATOM 1581 H21 LYS 235	37.080	17.978	1.622	1.00	0.00	BI
ATOM 1531 H ALA 230	35.871	27.026	7.654	1.00	0.00	BI	ATOM 1582 H22 LYS 235	36.854	16.466	2.363	1.00	0.00	BI
ATOM 1532 CA ALA 230	34.530	26.688	6.061	1.00	25.94	BI	ATOM 1583 H23 LYS 235	38.080	17.497	2.911	1.00	0.00	BI
ATOM 1533 CB ALA 230	35.193	27.852	5.312	1.00	15.76	BI	ATOM 1584 C LYS 235	30.363	18.847	3.204	1.00	45.20	BI
ATOM 1534 C ALA 230	34.794	25.403	5.304	1.00	29.42	BI	ATOM 1585 O LYS 235	29.722	18.102	2.463	1.00	45.60	BI
ATOM 1535 O ALA 230	34.014	25.061	4.423	1.00	32.07	BI	ATOM 1586 N LEU 236	27.807	19.332	4.301	1.00	33.54	BI
ATOM 1536 N ALA 231	35.878	24.671	5.572	1.00	32.16	BI	ATOM 1587 H LEU 236	30.363	19.888	4.885	1.00	0.00	BI
ATOM 1537 H ALA 231	36.556	25.045	6.175	1.00	0.00	BI	ATOM 1588 CA LEU 236	28.417	19.116	4.641	1.00	22.40	BI
ATOM 1538 CA ALA 231	36.141	23.364	4.957	1.00	31.99	BI	ATOM 1589 CB LEU 236	28.093	19.918	5.894	1.00	28.65	BI
ATOM 1539 CB ALA 231	37.489	22.847	5.428	1.00	32.77	BI	ATOM 1590 CG LEU 236	28.791	19.441	7.148	1.00	28.21	BI
ATOM 1540 C ALA 231	35.060	22.361	5.386	1.00	32.99	BI	ATOM 1591 CD1 LEU 236	28.703	20.460	8.268	1.00	24.14	BI
ATOM 1541 O ALA 231	34.599	21.575	4.576	1.00	34.12	BI	ATOM 1592 CD2 LEU 236	28.132	18.163	7.587	1.00	26.66	BI
ATOM 1542 N LEU 232	34.662	22.309	6.652	1.00	33.30	BI	ATOM 1593 C LEU 236	27.590	19.574	3.453	1.00	31.49	BI
ATOM 1543 H LEU 232	35.174	22.861	7.284	1.00	0.00	BI	ATOM 1594 O LEU 236	26.691	18.849	3.064	1.00	35.13	BI
ATOM 1544 CA LEU 232	33.558	21.506	7.165	1.00	35.33	BI	ATOM 1595 N CYS 237	27.870	20.670	2.753	1.00	34.49	BI
ATOM 1545 CB LEU 232	33.279	21.783	8.626	1.00	34.22	BI	ATOM 1596 H CYS 237	28.611	21.251	3.025	1.00	0.00	BI
ATOM 1546 CG LEU 232	32.410	20.861	9.394	1.00	33.16	BI	ATOM 1597 CA CYS 237	27.064	21.016	1.606	1.00	34.95	BI
ATOM 1547 CD1 LEU 232	33.191	19.545	9.451	1.00	34.59	BI	ATOM 1598 C CYS 237	27.324	20.090	0.451	1.00	35.97	BI
ATOM 1548 CD2 LEU 232	32.107	21.381	10.800	1.00	31.32	BI	ATOM 1599 O CYS 237	26.360	19.573	-0.089	1.00	36.09	BI
ATOM 1549 C LEU 232	32.271	21.829	6.440	1.00	36.65	BI	ATOM 1600 CB CYS 237	27.334	22.413	1.136	1.00	35.18	BI
ATOM 1550 O LEU 232	31.703	20.986	5.749	1.00	36.42	BI	ATOM 1601 SG CYS 237	26.409	22.880	-0.365	1.00	36.40	BI
ATOM 1551 N GLN 233	31.836	23.084	6.570	1.00	38.89	BI	ATOM 1602 N ALA 238	28.571	19.804	0.074	1.00	37.29	BI
ATOM 1552 H GLN 233	32.378	23.719	7.087	1.00	0.00	BI	ATOM 1603 H ALA 238	29.324	20.158	0.591	1.00	0.00	BI
ATOM 1553 CA GLN 233	30.637	23.579	5.933	1.00	40.02	BI	ATOM 1604 CA ALA 238	28.841	18.973	-1.090	1.00	36.80	BI
ATOM 1554 CB GLN 233	30.572	25.072	6.162	1.00	42.25	BI	ATOM 1605 CB ALA 238	30.274	18.684	-1.403	1.00	37.35	BI
ATOM 1555 CG GLN 233	30.290	25.398	7.626	1.00	48.22	BI	ATOM 1606 C ALA 238	28.320	17.617	-0.911	1.00	36.49	BI
ATOM 1556 CD GLN 233	30.021	26.879	7.983	1.00	53.75	BI	ATOM 1607 O ALA 238	27.645	17.198	-1.809	1.00	36.54	BI
ATOM 1557 OE1 GLN 233	30.799	27.810	7.718	1.00	55.93	BI	ATOM 1608 N THR 239	28.628	16.969	0.193	1.00	38.80	BI
ATOM 1558 NE2 GLN 233	28.909	27.215	8.634	1.00	56.51	BI	ATOM 1609 H THR 239	29.236	17.391	0.821	1.00	0.00	BI
ATOM 1559 HE1 GLN 233	28.810	28.144	8.902	1.00	0.00	BI	ATOM 1610 CA THR 239	28.230	15.587	0.464	1.00	41.31	BI
ATOM 1560 HE2 GLN 233	28.705	26.533	8.710	1.00	0.00	BI	ATOM 1611 CB THR 239	29.158	15.035	1.554	1.00	42.38	BI
ATOM 1561 C GLN 233	30.635	23.243	4.441	1.00	39.70	BI	ATOM 1612 OG1 THR 239	30.473	15.265	1.031	1.00	45.70	BI
ATOM 1562 O GLN 233	29.631	22.777	3.898	1.00	40.20	BI	ATOM 1613 HG1 THR 239	31.019	15.668	1.709	1.00	0.00	BI
ATOM 1563 N GLU 234	31.744	23.377	3.736	1.00	39.32	BI	ATOM 1614 CG2 THR 239	28.936	13.574	1.916	1.00	41.85	BI
ATOM 1564 H GLU 234	32.544	23.750	4.163	1.00	0.00	BI	ATOM 1615 C THR 239	26.771	15.341	0.864	1.00	41.94	BI
ATOM 1565 CA GLU 234	31.809	23.025	2.329	1.00	39.23	BI	ATOM 1616 O THR 239	26.260	14.284	0.460	1.00	43.34	BI
ATOM 1566 CB GLU 234	33.155	23.434	1.811	1.00	40.25	BI	ATOM 1617 N TYR 240	26.095	16.207	1.669	1.00	40.07	BI
ATOM 1567 CG GLU 234	33.292	23.028	0.383	1.00	47.69	BI	ATOM 1618 H TYR 240	26.538	17.034	1.953	1.00	0.00	BI
ATOM 1568 CD GLU 234	34.733	23.056	-0.073	1.00	53.40	BI	ATOM 1619 CA TYR 240	24.718	15.992	2.084	1.00	38.21	BI
ATOM 1569 OE1 GLU 234	34.986	23.721	-1.100	1.00	53.78	BI	ATOM 1620 CB TYR 240	24.594	15.993	3.618	1.00	38.08	BI
ATOM 1570 OE2 GLU 234	35.568	22.400	0.590	1.00	57.55	BI	ATOM 1621 CG TYR 240	25.524	14.926	4.193	1.00	43.37	BI
ATOM 1571 C GLU 234	31.580	21.535	2.136	1.00	37.09	BI	ATOM 1622 CD1 TYR 240	26.475	15.243	5.149	1.00	45.06	BI
ATOM 1572 O GLU 234	30.884	21.217	1.188	1.00	36.67	BI	ATOM 1623 CE1 TYR 240	27.470	14.283	5.529	1.00	47.35	BI
ATOM 1573 N LYS 235	32.092	20.623	2.986	1.00	37.27	BI	ATOM 1624 CD2 TYR 240	25.518	13.643	3.641	1.00	45.89	BI
ATOM 1574 H LYS 235	32.668	20.965	3.706	1.00	0.00	BI	ATOM 1625 CE2 TYR 240	26.442	12.690	4.003	1.00	44.77	BI
ATOM 1575 CA LYS 235	31.832	19.177	2.942	1.00	36.27	BI	ATOM 1626 CE3 TYR 240	27.410	13.005	4.543	1.00	47.96	BI
ATOM 1576 CB LYS 235	32.516	18.365	3.997	1.00	34.92	BI	ATOM 1627 OH TYR 240	28.390	12.047	5.244	1.00	46.51	BI
ATOM 1577 CG LYS 235	33.978	18.483	4.107	1.00	38.47	BI	ATOM 1628 HII TYR 240	28.027	11.187	4.992	1.00	0.00	BI
ATOM 1578 CD LYS 235	34.762	17.999	2.921	1.00	38.07	BI	ATOM 1629 C TYR 240	23.781	17.632	1.516	1.00	39.49	BI
ATOM 1579 CE LYS 235	36.192	18.051	3.460	1.00	39.15	BI	ATOM 1630 O TYR 240	22.587	16.934	1.775	1.00	42.76	BI

FIGURE 5

ATOM 1733 C LEU 251	17.104	22.372	15.493	1.00	42.78	BI	ATOM 1784 N PRO 258	23.441	24.392	22.608	1.00	43.05	BI
ATOM 1734 O LEU 251	17.124	21.554	16.395	1.00	45.44	BI	ATOM 1785 CD PRO 258	24.133	23.371	23.296	1.00	43.29	BI
ATOM 1735 N GLY 252	17.826	23.477	15.610	1.00	44.86	BI	ATOM 1786 CA PRO 258	23.559	25.616	23.360	1.00	43.82	BI
ATOM 1736 H GLY 252	17.750	24.160	14.910	1.00	0.00	BI	ATOM 1787 CB PRO 258	24.295	25.736	24.612	1.00	41.97	BI
ATOM 1737 CA GLY 252	18.734	23.711	16.719	1.00	46.68	BI	ATOM 1788 CG PRO 258	25.107	24.064	24.186	1.00	42.71	BI
ATOM 1738 C GLY 252	18.071	23.596	18.067	1.00	49.18	BI	ATOM 1789 C PRO 258	24.527	26.703	22.555	1.00	46.06	BI
ATOM 1739 O GLY 252	18.709	23.318	19.077	1.00	49.23	BI	ATOM 1790 O PRO 258	24.983	26.513	21.560	1.00	46.59	BI
ATOM 1740 N HIS 253	16.756	23.787	18.046	1.00	53.74	BI	ATOM 1791 N TRP 259	23.996	27.887	23.106	1.00	46.75	BI
ATOM 1741 H HIS 253	16.358	24.055	17.190	1.00	0.00	BI	ATOM 1792 H TRP 259	23.588	27.921	23.994	1.00	41.00	BI
ATOM 1742 CA HIS 253	15.859	23.649	19.197	1.00	57.46	BI	ATOM 1793 CA TRP 259	24.427	29.143	22.517	1.00	45.77	BI
ATOM 1743 CB HIS 253	14.468	24.157	18.764	1.00	62.93	BI	ATOM 1794 CB TRP 259	23.213	30.071	22.397	1.00	46.60	BI
ATOM 1744 CG HIS 253	13.212	23.813	19.577	1.00	68.75	BI	ATOM 1795 CG TRP 259	23.556	31.372	21.749	1.00	47.51	BI
ATOM 1745 CD HIS 253	12.031	24.529	19.414	1.00	71.00	BI	ATOM 1796 CD2 TRP 259	23.860	31.525	20.430	1.00	47.83	BI
ATOM 1746 ND1 HIS 253	12.980	22.854	20.479	1.00	70.67	BI	ATOM 1797 CE2 TRP 259	24.154	32.888	20.392	1.00	48.47	BI
ATOM 1747 HD1 HIS 253	13.627	22.193	20.830	1.00	0.00	BI	ATOM 1798 CE3 TRP 259	23.940	30.745	19.290	1.00	47.59	BI
ATOM 1748 CE1 HIS 253	11.723	22.966	20.845	1.00	73.40	BI	ATOM 1799 CD1 TRP 259	23.639	32.520	22.493	1.00	48.64	BI
ATOM 1749 NE1 HIS 253	11.156	23.973	20.204	1.00	72.91	BI	ATOM 1800 NE1 TRP 259	24.013	33.421	21.628	1.00	48.27	BI
ATOM 1750 HE2 HIS 253	10.218	24.260	20.311	1.00	0.00	BI	ATOM 1801 HE1 TRP 259	24.224	34.344	21.870	1.00	0.00	BI
ATOM 1751 C HIS 253	15.771	22.209	19.691	1.00	56.06	BI	ATOM 1802 CZ2 TRP 259	24.531	33.486	19.195	1.00	47.40	BI
ATOM 1752 O HIS 253	15.880	21.827	20.857	1.00	56.17	BI	ATOM 1803 CZ3 TRP 259	24.317	31.344	18.097	1.00	49.07	BI
ATOM 1753 N SER 254	15.395	21.435	18.724	1.00	53.46	BI	ATOM 1804 C112 TRP 259	24.613	32.706	18.050	1.00	49.12	BI
ATOM 1754 H SER 254	15.278	21.783	17.813	1.00	0.00	BI	ATOM 1805 C TRP 259	25.459	29.727	23.440	1.00	44.01	BI
ATOM 1755 CA SER 254	15.177	20.034	18.898	1.00	52.61	BI	ATOM 1806 O TRP 259	25.340	29.664	24.671	1.00	43.25	BI
ATOM 1756 CB SER 254	14.613	19.595	17.576	1.00	53.04	BI	ATOM 1807 N ALA 260	26.469	30.247	22.777	1.00	43.01	BI
ATOM 1757 CG SER 254	13.794	20.686	17.158	1.00	56.04	BI	ATOM 1808 H ALA 260	26.523	30.198	21.796	1.00	0.00	BI
ATOM 1758 HG SER 254	13.369	20.467	16.319	1.00	0.00	BI	ATOM 1809 CA ALA 260	27.493	30.973	23.482	1.00	43.48	BI
ATOM 1759 C SER 254	16.312	19.386	19.275	1.00	51.48	BI	ATOM 1810 CB ALA 260	28.874	30.549	22.969	1.00	43.31	BI
ATOM 1760 O SER 254	16.596	18.639	20.245	1.00	51.90	BI	ATOM 1811 C ALA 260	27.249	32.486	23.216	1.00	43.41	BI
ATOM 1761 N LEU 255	17.577	19.790	18.562	1.00	49.31	BI	ATOM 1812 O ALA 260	27.315	32.946	22.054	1.00	40.55	BI
ATOM 1762 H LEU 255	17.430	20.480	17.889	1.00	0.00	BI	ATOM 1813 N PRO 261	26.853	33.267	24.253	1.00	42.61	BI
ATOM 1763 CA LEU 255	18.913	19.272	18.723	1.00	46.02	BI	ATOM 1814 CD PRO 261	26.527	32.807	25.606	1.00	42.33	BI
ATOM 1764 CB LEU 255	19.706	19.723	17.537	1.00	44.66	BI	ATOM 1815 CA PRO 261	26.720	34.701	24.199	1.00	42.37	BI
ATOM 1765 CG LEU 255	19.362	18.968	16.274	1.00	44.51	BI	ATOM 1816 CB PRO 261	25.778	34.987	25.335	1.00	41.46	BI
ATOM 1766 CD1 LEU 255	19.810	19.679	15.006	1.00	43.16	BI	ATOM 1817 CG PRO 261	26.251	34.060	26.411	1.00	40.00	BI
ATOM 1767 CD2 LEU 255	19.969	17.604	16.456	1.00	44.67	BI	ATOM 1818 C PRO 261	28.087	35.369	24.311	1.00	42.22	BI
ATOM 1768 C LEU 255	19.536	19.718	20.012	1.00	46.56	BI	ATOM 1819 O PRO 261	28.988	34.956	25.037	1.00	38.82	BI
ATOM 1769 O LEU 255	20.365	19.174	20.440	1.00	46.82	BI	ATOM 1820 N LEU 262	28.234	36.403	23.486	1.00	45.20	BI
ATOM 1770 N GLY 256	18.918	20.259	20.581	1.00	45.93	BI	ATOM 1821 H LEU 262	27.513	36.610	22.853	1.00	0.00	BI
ATOM 1771 H GLY 256	18.210	21.225	20.101	1.00	0.00	BI	ATOM 1822 CA LEU 262	29.434	37.210	23.498	1.00	46.50	BI
ATOM 1772 CA GLY 256	19.277	21.273	21.890	1.00	46.68	BI	ATOM 1823 CB LEU 262	30.531	36.609	22.610	1.00	45.09	BI
ATOM 1773 C GLY 256	20.669	21.866	21.970	1.00	47.28	BI	ATOM 1824 CG LEU 262	31.903	37.157	22.964	1.00	42.55	BI
ATOM 1774 O GLY 256	21.273	21.844	23.056	1.00	49.64	BI	ATOM 1825 CD1 LEU 262	32.344	36.695	24.338	1.00	41.52	BI
ATOM 1775 N ILE 257	21.143	22.441	20.849	1.00	45.74	BI	ATOM 1826 CD2 LEU 262	32.850	36.730	21.900	1.00	44.21	BI
ATOM 1776 H ILE 257	20.497	22.589	20.128	1.00	0.00	BI	ATOM 1827 C LEU 262	29.154	38.628	23.035	1.00	48.56	BI
ATOM 1777 CA ILE 257	22.481	23.017	20.776	1.00	43.64	BI	ATOM 1828 O LEU 262	29.633	39.470	23.790	1.00	48.23	BI
ATOM 1778 CB ILE 257	22.684	23.363	19.257	1.00	42.54	BI	ATOM 1829 N SER 263	28.388	38.956	21.960	1.00	51.33	BI
ATOM 1779 CG1 ILE 257	23.988	24.110	19.073	1.00	41.05	BI	ATOM 1830 H SER 263	27.982	38.242	21.427	1.00	0.00	BI
ATOM 1780 CG1 ILE 257	22.694	22.088	18.437	1.00	40.55	BI	ATOM 1831 CA SER 263	28.127	40.339	21.494	1.00	55.19	BI
ATOM 1781 CD ILE 257	22.452	22.468	16.970	1.00	39.49	BI	ATOM 1832 CB SER 263	26.871	40.511	20.621	1.00	57.17	BI
ATOM 1782 C ILE 257	22.559	24.246	21.616	1.00	43.27	BI	ATOM 1833 CG SER 263	26.498	39.411	19.776	1.00	64.12	BI
ATOM 1783 O ILE 257	21.706	25.110	21.450	1.00	43.22	BI	ATOM 1834 HG SER 263	26.093	38.741	20.336	1.00	0.00	BI

FIGURE 5

ATOM 1835 C SER 263	27.909	41.354	22.600	1.00	56.15	B1	ATOM 1886 C LEU 276	37.673	35.833	28.638	1.00	47.84	B2
ATOM 1836 O SER 263	28.744	42.243	22.753	1.00	57.88	B1	ATOM 1887 O LEU 276	37.784	34.803	27.964	1.00	48.51	B2
ATOM 1837 N SER 264	26.899	41.231	23.452	1.00	56.52	B1	ATOM 1888 N ALA 277	37.074	35.840	29.804	1.00	45.56	B2
ATOM 1838 H SER 264	26.277	40.478	23.415	1.00	0.00	B1	ATOM 1889 H ALA 277	36.898	36.662	30.389	1.00	0.00	B2
ATOM 1839 CA SER 264	26.716	42.204	24.494	1.00	58.28	B1	ATOM 1890 CA ALA 277	36.613	34.605	30.365	1.00	45.77	B2
ATOM 1840 CB SER 264	25.313	41.977	25.064	1.00	58.77	B1	ATOM 1891 CB ALA 277	36.147	34.810	31.783	1.00	47.87	B2
ATOM 1841 CG SER 264	25.099	40.726	25.713	1.00	58.50	B1	ATOM 1892 C ALA 277	35.442	34.111	29.542	1.00	45.33	B2
ATOM 1842 HG SER 264	25.385	40.832	26.632	1.00	0.00	B1	ATOM 1893 O ALA 277	35.342	32.926	29.271	1.00	44.20	B2
ATOM 1843 C SER 264	27.800	42.168	25.584	1.00	59.95	B1	ATOM 1894 N GIN 278	34.592	35.000	29.049	1.00	45.14	B2
ATOM 1844 O SER 264	27.610	42.805	26.620	1.00	60.44	B1	ATOM 1895 H GIN 278	34.231	35.944	29.263	1.00	0.00	B2
ATOM 1845 N CYS 265	28.948	41.484	25.466	1.00	61.37	B1	ATOM 1896 CA GIN 278	33.435	34.601	28.284	1.00	45.27	B2
ATOM 1846 H CYS 265	29.192	41.114	24.596	1.00	0.00	B1	ATOM 1897 CB GIN 278	32.550	35.825	28.083	1.00	48.13	B2
ATOM 1847 CA CYS 265	29.958	41.502	26.509	1.00	62.57	B1	ATOM 1898 CG GIN 278	31.140	35.442	28.484	1.00	56.00	B2
ATOM 1848 CB CYS 265	30.991	40.418	26.285	1.00	64.32	B1	ATOM 1899 CD GIN 278	30.045	36.464	28.178	1.00	61.94	B2
ATOM 1849 SG CYS 265	32.322	40.638	27.504	1.00	71.40	B1	ATOM 1900 OEI GIN 278	29.048	36.530	28.896	1.00	65.95	B2
ATOM 1850 C CYS 265	30.667	42.860	26.515	1.00	63.12	B1	ATOM 1901 NEI GIN 278	30.080	37.291	27.132	1.00	65.55	B2
ATOM 1851 OTI CYS 265	31.065	43.360	25.444	1.00	63.44	B1	ATOM 1902 HEI2 GIN 278	30.829	37.221	26.510	1.00	0.00	B2
ATOM 1852 OT2 CYS 265	30.809	43.408	27.610	1.00	61.72	B1	ATOM 1903 HEI2 GIN 278	29.343	37.927	27.056	1.00	0.00	B2
ATOM 1853 CB ALA 272	40.070	43.327	30.788	1.00	77.44	B2	ATOM 1904 C GIN 278	33.812	33.971	26.950	1.00	43.16	B2
ATOM 1854 CA 272	38.698	41.201	30.601	1.00	76.53	B2	ATOM 1905 O GIN 278	33.173	33.050	26.462	1.00	40.58	B2
ATOM 1855 O ALA 272	37.525	40.873	30.361	1.00	76.81	B2	ATOM 1906 N LEU 279	34.869	34.476	26.331	1.00	43.32	B2
ATOM 1856 HTI ALA 272	37.486	41.550	30.261	1.00	0.00	B2	ATOM 1907 H LEU 279	35.328	35.227	26.767	1.00	0.00	B2
ATOM 1857 HT2 ALA 272	37.357	42.450	28.996	1.00	0.00	B2	ATOM 1908 CA LEU 279	35.398	33.966	25.069	1.00	42.80	B2
ATOM 1858 N ALA 272	37.973	43.169	29.427	1.00	76.81	B2	ATOM 1909 CB LEU 279	36.363	34.790	24.626	1.00	41.42	B2
ATOM 1859 HT3 ALA 272	38.195	43.924	28.752	1.00	0.00	B2	ATOM 1910 CG LEU 279	36.885	35.014	23.190	1.00	40.76	B2
ATOM 1860 CA ALA 272	39.176	42.460	29.853	1.00	77.02	B2	ATOM 1911 CDI LEU 279	38.239	35.647	23.130	1.00	41.76	B2
ATOM 1861 N ALA 273	39.485	40.547	31.487	1.00	74.93	B2	ATOM 1912 CD2 LEU 279	36.943	33.753	22.411	1.00	40.01	B2
ATOM 1862 H ALA 273	40.334	40.963	31.745	1.00	0.00	B2	ATOM 1913 C LEU 279	35.876	32.554	25.341	1.00	42.92	B2
ATOM 1863 CA ALA 273	39.244	39.241	32.119	1.00	72.64	B2	ATOM 1914 O LEU 279	35.572	31.598	24.640	1.00	42.57	B2
ATOM 1864 CB ALA 273	39.704	39.279	33.558	1.00	71.92	B2	ATOM 1915 N HIS 280	36.654	32.463	26.403	1.00	43.93	B2
ATOM 1865 C ALA 273	37.872	38.599	32.118	1.00	71.60	B2	ATOM 1916 H HIS 280	36.837	33.282	26.917	1.00	0.00	B2
ATOM 1866 O ALA 273	37.806	37.458	31.702	1.00	71.68	B2	ATOM 1917 CA HIS 280	37.215	31.223	26.850	1.00	46.12	B2
ATOM 1867 N GLY 274	36.775	39.282	32.484	1.00	70.20	B2	ATOM 1918 CB HIS 280	38.029	31.506	28.101	1.00	48.74	B2
ATOM 1868 H GLY 274	36.903	40.167	32.874	1.00	0.00	B2	ATOM 1919 CG HIS 280	38.914	30.320	28.394	1.00	54.16	B2
ATOM 1869 CA GLY 274	35.412	38.758	32.425	1.00	66.78	B2	ATOM 1920 CD2 HIS 280	40.041	30.069	27.650	1.00	56.02	B2
ATOM 1870 C GLY 274	35.050	38.437	30.990	1.00	65.05	B2	ATOM 1921 NDI HIS 280	38.759	29.326	29.264	1.00	56.01	B2
ATOM 1871 O GLY 274	34.627	37.320	30.709	1.00	66.44	B2	ATOM 1922 HD1 HIS 280	38.012	29.203	29.890	1.00	0.00	B2
ATOM 1872 N CYS 275	35.301	39.364	30.048	1.00	62.77	B2	ATOM 1923 CEI HIS 280	39.744	28.483	29.058	1.00	56.64	B2
ATOM 1873 H CYS 275	35.634	40.223	30.357	1.00	0.00	B2	ATOM 1924 NEI HIS 280	40.507	28.937	28.088	1.00	56.64	B2
ATOM 1874 CA CYS 275	35.026	39.188	28.611	1.00	59.30	B2	ATOM 1925 HEI2 HIS 280	41.282	28.478	27.684	1.00	0.00	B2
ATOM 1875 C CYS 275	35.875	38.063	28.054	1.00	55.89	B2	ATOM 1926 C HIS 280	36.161	30.134	27.117	1.00	45.65	B2
ATOM 1876 O CYS 275	35.425	37.152	27.351	1.00	54.41	B2	ATOM 1927 O HIS 280	36.362	28.977	26.711	1.00	46.23	B2
ATOM 1877 CB CYS 275	35.349	40.466	27.827	1.00	61.50	B2	ATOM 1928 N SER 281	35.086	30.473	27.822	1.00	43.91	B2
ATOM 1878 SG CYS 275	34.119	40.937	26.577	1.00	66.63	B2	ATOM 1929 H SER 281	33.009	31.467	28.219	1.00	0.00	B2
ATOM 1879 N LEU 276	37.124	38.114	28.506	1.00	52.23	B2	ATOM 1930 CA SER 281	34.008	29.574	28.105	1.00	41.53	B2
ATOM 1880 H LEU 276	37.350	38.722	29.233	1.00	0.00	B2	ATOM 1931 CB SER 281	35.026	30.291	29.002	1.00	44.18	B2
ATOM 1881 CA LEU 276	38.091	37.163	28.066	1.00	48.93	B2	ATOM 1932 CG SER 281	33.761	30.812	30.113	1.00	47.74	B2
ATOM 1882 CB LEU 276	39.483	37.564	28.542	1.00	45.96	B2	ATOM 1933 IIG SER 281	33.288	30.648	30.931	1.00	0.00	B2
ATOM 1883 CG LEU 276	40.241	38.557	27.670	1.00	43.20	B2	ATOM 1934 C SER 281	33.382	29.169	26.787	1.00	43.35	B2
ATOM 1884 CYN I1U 276	41.599	38.782	28.279	1.00	44.63	B2	ATOM 1935 O SER 281	33.334	27.973	26.496	1.00	44.83	B2
ATOM 1885 C12 I1U 276	40.429	38.033	26.271	1.00	40.55	B2	ATOM 1936 N GLY 282	37.977	30.120	25.940	1.00	42.53	B2

FIGURE 5

ATOM 1937 H GLY 282	33.043	31.058	26.221	1.00	0.00	B2
ATOM 1938 CA GLY 282	32.363	29.869	24.632	1.00	40.65	B2
ATOM 1939 C GLY 282	33.175	28.937	23.755	1.00	39.06	B2
ATOM 1940 O GLY 282	32.584	28.075	23.107	1.00	40.10	B2
ATOM 1941 N LEU 283	34.514	29.066	23.776	1.00	37.39	B2
ATOM 1942 H LEU 283	34.880	29.807	24.304	1.00	0.00	B2
ATOM 1943 CA LEU 283	35.465	28.213	23.037	1.00	35.06	B2
ATOM 1944 CB LEU 283	36.902	28.718	23.089	1.00	30.20	B2
ATOM 1945 CG LEU 283	37.167	30.001	22.302	1.00	25.73	B2
ATOM 1946 CD1 LEU 283	38.539	30.461	22.664	1.00	24.38	B2
ATOM 1947 CD2 LEU 283	37.036	29.402	20.815	1.00	21.94	B2
ATOM 1948 C LEU 283	35.470	26.851	23.651	1.00	34.81	B2
ATOM 1949 O LEU 283	35.314	25.859	22.947	1.00	31.09	B2
ATOM 1950 N PHE 284	35.533	26.842	24.973	1.00	37.62	B2
ATOM 1951 H PHE 284	35.567	27.686	25.467	1.00	0.00	B2
ATOM 1952 CA PHE 284	35.485	25.596	25.710	1.00	42.51	B2
ATOM 1953 CB PHE 284	35.542	25.877	27.184	1.00	49.49	B2
ATOM 1954 CG PHE 284	36.221	24.770	27.968	1.00	58.39	B2
ATOM 1955 CD1 PHE 284	37.265	25.108	28.816	1.00	63.05	B2
ATOM 1956 CD2 PHE 284	35.810	23.453	27.861	1.00	60.84	B2
ATOM 1957 CE1 PHE 284	37.900	24.124	29.563	1.00	65.86	B2
ATOM 1958 CE2 PHE 284	36.444	22.480	28.605	1.00	64.49	B2
ATOM 1959 CZ PHE 284	37.486	22.810	29.455	1.00	66.32	B2
ATOM 1960 C PHE 284	34.204	24.849	25.384	1.00	41.44	B2
ATOM 1961 O PHE 284	34.257	23.630	25.306	1.00	41.42	B2
ATOM 1962 N LEU 285	33.100	25.563	25.101	1.00	41.24	B2
ATOM 1963 H LEU 285	33.192	26.534	25.174	1.00	0.00	B2
ATOM 1964 CA LEU 285	31.781	25.025	24.730	1.00	38.92	B2
ATOM 1965 CB LEU 285	30.727	26.139	24.807	1.00	39.05	B2
ATOM 1966 CG LEU 285	29.292	25.740	24.481	1.00	41.16	B2
ATOM 1967 CD1 LEU 285	28.711	24.981	25.662	1.00	41.12	B2
ATOM 1968 CD2 LEU 285	28.472	26.971	24.139	1.00	39.60	B2
ATOM 1969 C LEU 285	31.780	24.441	23.379	1.00	37.34	B2
ATOM 1970 O LEU 285	31.245	23.351	23.095	1.00	36.97	B2
ATOM 1971 N TYR 286	32.352	25.172	22.372	1.00	35.26	B2
ATOM 1972 H TYR 286	32.705	26.062	22.593	1.00	0.00	B2
ATOM 1973 CA TYR 286	32.455	24.660	21.033	1.00	35.04	B2
ATOM 1974 CB TYR 286	32.891	25.790	20.122	1.00	34.44	B2
ATOM 1975 CG TYR 286	31.690	26.684	19.008	1.00	34.75	B2
ATOM 1976 CD1 TYR 286	31.433	27.879	20.469	1.00	35.67	B2
ATOM 1977 CE1 TYR 286	30.313	28.620	20.158	1.00	36.90	B2
ATOM 1978 CD2 TYR 286	30.823	26.255	18.839	1.00	36.19	B2
ATOM 1979 CE2 TYR 286	29.707	26.990	18.521	1.00	37.55	B2
ATOM 1980 CZ TYR 286	29.449	28.164	19.178	1.00	37.73	B2
ATOM 1981 OH TYR 286	28.285	28.826	18.823	1.00	38.04	B2
ATOM 1982 HH TYR 286	28.289	29.707	19.243	1.00	0.00	B2
ATOM 1983 C TYR 286	33.393	23.464	20.976	1.00	34.80	B2
ATOM 1984 O TYR 286	33.071	22.537	20.180	1.00	35.35	B2
ATOM 1985 N ALA 287	34.527	23.339	21.636	1.00	34.66	B2
ATOM 1986 H ALA 287	34.803	24.088	22.206	1.00	0.00	B2
ATOM 1987 CA ALA 287	35.350	22.108	21.565	1.00	34.28	B2
ATOM 1988 CB ALA 287	36.617	22.291	22.415	1.00	33.63	B2
ATOM 1989 C ALA 287	34.528	20.906	22.073	1.00	34.32	B2
ATOM 1990 O ALA 287	34.535	19.827	21.478	1.00	33.07	B2
ATOM 1991 N GLY 288	33.723	21.118	23.111	1.00	33.19	B2
ATOM 1992 H GLY 288	33.791	21.985	23.564	1.00	0.00	B2
ATOM 1993 CA GLY 288	32.761	20.162	23.655	1.00	35.62	B2
ATOM 1994 C GLY 288	31.744	19.606	22.636	1.00	36.81	B2
ATOM 1995 O GLY 288	31.037	18.379	22.444	1.00	34.97	B2
ATOM 1996 N LEU 289	31.037	20.536	21.966	1.00	36.69	B2
ATOM 1997 H LEU 289	31.200	21.476	22.201	1.00	0.00	B2
ATOM 1998 CA LEU 289	30.018	20.249	20.954	1.00	35.05	B2
ATOM 1999 CB LEU 289	29.351	21.576	20.502	1.00	36.32	B2
ATOM 2000 CG LEU 289	28.552	22.450	21.464	1.00	35.76	B2
ATOM 2001 CD1 LEU 289	28.256	23.821	20.890	1.00	32.66	B2
ATOM 2002 CD2 LEU 289	27.246	21.780	21.697	1.00	35.35	B2
ATOM 2003 C LEU 289	30.536	19.519	19.714	1.00	34.21	B2
ATOM 2004 O LEU 289	29.871	18.694	19.078	1.00	33.28	B2
ATOM 2005 N LEU 290	31.756	19.902	19.355	1.00	33.25	B2
ATOM 2006 H LEU 290	32.183	20.634	19.850	1.00	0.00	B2
ATOM 2007 CA LEU 290	32.448	19.345	18.230	1.00	32.44	B2
ATOM 2008 CB LEU 290	33.779	20.159	18.000	1.00	32.62	B2
ATOM 2009 CG LEU 290	33.560	21.509	17.315	1.00	32.05	B2
ATOM 2010 CD1 LEU 290	34.889	22.189	17.349	1.00	32.58	B2
ATOM 2011 CD2 LEU 290	33.068	21.374	15.879	1.00	31.74	B2
ATOM 2012 C LEU 290	32.737	17.908	18.558	1.00	31.94	B2
ATOM 2013 O LEU 290	32.432	17.020	17.772	1.00	30.50	B2
ATOM 2014 N GLN 291	33.249	17.711	19.770	1.00	33.58	B2
ATOM 2015 H GLN 291	33.512	18.494	20.298	1.00	0.00	B2
ATOM 2016 CA GLN 291	33.499	16.372	20.311	1.00	36.39	B2
ATOM 2017 CB GLN 291	33.988	16.490	21.702	1.00	36.86	B2
ATOM 2018 CG GLN 291	34.926	15.367	21.950	1.00	39.48	B2
ATOM 2019 CD GLN 291	35.658	15.503	23.252	1.00	40.79	B2
ATOM 2020 OE1 GLN 291	36.457	14.626	23.549	1.00	44.80	B2
ATOM 2021 NE1 GLN 291	35.494	16.535	24.072	1.00	42.59	B2
ATOM 2022 HE1 GLN 291	34.928	17.287	23.817	1.00	0.00	B2
ATOM 2023 HE2 GLN 291	35.910	16.463	24.958	1.00	0.00	B2
ATOM 2024 C GLN 291	32.233	15.536	20.307	1.00	36.66	B2
ATOM 2025 O GLN 291	32.220	14.478	19.707	1.00	37.46	B2
ATOM 2026 N ALA 292	31.143	16.023	20.913	1.00	38.37	B2
ATOM 2027 H ALA 292	31.255	16.849	21.418	1.00	0.00	B2
ATOM 2028 CA ALA 292	29.778	15.451	20.857	1.00	39.25	B2
ATOM 2029 CB ALA 292	28.818	16.483	21.444	1.00	40.28	B2
ATOM 2030 C ALA 292	29.215	14.999	19.484	1.00	38.65	B2
ATOM 2031 O ALA 292	28.411	14.067	19.356	1.00	37.58	B2
ATOM 2032 H LEU 293	29.614	15.702	18.430	1.00	39.00	B2
ATOM 2033 N LEU 293	30.149	16.513	18.574	1.00	0.00	B2
ATOM 2034 CA LEU 293	29.265	15.335	17.077	1.00	39.74	B2
ATOM 2035 CB LEU 293	29.662	16.418	16.106	1.00	37.53	B2
ATOM 2036 CG LEU 293	28.969	17.701	16.138	1.00	34.34	B2
ATOM 2037 CD1 LEU 293	29.547	18.582	15.053	1.00	33.88	B2
ATOM 2038 CD2 LEU 293	27.503	17.462	15.918	1.00	35.65	B2

FIGURE 5

ATOM 2039 C LEU 293	29.933	14.060	16.596	1.00	-40.86	B2	ATOM 2090 C GLN 299	34.923	12.453	6.160	1.00	51.04	B2
ATOM 2040 O LEU 293	29.686	13.669	15.449	1.00	40.58	B2	ATOM 2091 O GLN 299	35.796	13.186	5.718	1.00	53.38	B2
ATOM 2041 N GLU 294	30.887	13.495	17.365	1.00	42.12	B2	ATOM 2092 N LEU 300	34.118	12.918	7.120	1.00	48.45	B2
ATOM 2042 H GLU 294	31.131	13.963	18.190	1.00	0.00	B2	ATOM 2093 H LEU 300	33.382	12.351	7.437	1.00	0.00	B2
ATOM 2043 CA GLU 294	31.598	12.253	17.076	1.00	42.89	B2	ATOM 2094 CA LEU 300	34.272	14.220	7.745	1.00	43.32	B2
ATOM 2044 CB GLU 294	30.806	10.984	17.485	1.00	46.38	B2	ATOM 2095 CB LEU 300	32.856	14.719	8.021	1.00	41.39	B2
ATOM 2045 CG GLU 294	30.715	10.614	18.972	1.00	56.76	B2	ATOM 2096 CG LEU 300	32.073	15.546	6.974	1.00	37.99	B2
ATOM 2046 CD GLU 294	29.271	10.408	19.486	1.00	63.70	B2	ATOM 2097 CD LEU 300	31.872	14.824	5.688	1.00	48.34	B2
ATOM 2047 OEI GLU 294	29.058	10.603	20.702	1.00	67.72	B2	ATOM 2098 OEI LEU 300	30.705	15.809	7.522	1.00	37.67	B2
ATOM 2048 OEI GLU 294	28.363	10.074	18.692	1.00	64.81	B2	ATOM 2099 C LEU 300	35.142	14.220	9.019	1.00	42.84	B2
ATOM 2049 C GLU 294	31.972	12.068	15.632	1.00	41.53	B2	ATOM 2100 O LEU 300	35.558	15.278	9.541	1.00	41.56	B2
ATOM 2050 O GLU 294	31.804	11.007	15.021	1.00	40.29	B2	ATOM 2101 N GLY 301	35.467	13.016	9.578	1.00	40.83	B2
ATOM 2051 N GLY 295	32.424	13.203	15.106	1.00	40.93	B2	ATOM 2102 H GLY 301	35.157	12.221	9.046	1.00	0.00	B2
ATOM 2052 H GLY 295	32.357	14.033	15.621	1.00	0.00	B2	ATOM 2103 CA GLY 301	36.199	12.826	10.779	1.00	36.72	B2
ATOM 2053 CA GLY 295	32.998	13.236	13.783	1.00	39.95	B2	ATOM 2104 C GLY 301	37.500	13.607	10.887	1.00	37.69	B2
ATOM 2054 C GLY 295	32.027	13.230	12.634	1.00	40.60	B2	ATOM 2105 O GLY 301	37.665	14.406	11.809	1.00	37.31	B2
ATOM 2055 O GLY 295	32.477	13.216	11.487	1.00	40.96	B2	ATOM 2106 N PRO 302	38.468	13.452	9.985	1.00	37.33	B2
ATOM 2056 N ILE 296	30.728	13.296	12.898	1.00	41.18	B2	ATOM 2107 CD PRO 302	38.353	12.630	8.790	1.00	37.77	B2
ATOM 2057 H ILE 296	30.446	13.210	13.825	1.00	0.00	B2	ATOM 2108 CA PRO 302	39.676	14.281	9.864	1.00	37.60	B2
ATOM 2058 CA ILE 296	29.687	13.306	11.888	1.00	44.02	B2	ATOM 2109 CB PRO 302	40.256	13.907	8.541	1.00	36.62	B2
ATOM 2059 CB ILE 296	29.683	14.580	11.009	1.00	43.49	B2	ATOM 2110 CG PRO 302	39.047	13.487	7.745	1.00	37.94	B2
ATOM 2060 CG2 ILE 296	28.288	14.685	10.421	1.00	40.56	B2	ATOM 2111 C PRO 302	39.486	15.782	10.033	1.00	37.45	B2
ATOM 2061 CG1 ILE 296	30.047	15.831	11.793	1.00	45.11	B2	ATOM 2112 O PRO 302	40.132	16.398	10.901	1.00	38.33	B2
ATOM 2062 CD ILE 296	30.039	17.189	11.062	1.00	46.06	B2	ATOM 2113 N THR 303	38.547	16.311	9.204	1.00	37.05	B2
ATOM 2063 C ILE 296	29.820	12.107	10.949	1.00	46.71	B2	ATOM 2114 H THR 303	38.085	15.727	8.567	1.00	0.00	B2
ATOM 2064 O ILE 296	28.918	11.279	11.060	1.00	50.61	B2	ATOM 2115 CA THR 303	38.119	17.705	9.128	1.00	35.81	B2
ATOM 2065 N SER 297	30.767	11.475	10.019	1.00	47.21	B2	ATOM 2116 CB THR 303	36.963	17.770	8.123	1.00	37.26	B2
ATOM 2066 H SER 297	31.526	12.491	9.936	1.00	0.00	B2	ATOM 2117 CG1 THR 303	37.416	17.161	6.909	1.00	39.64	B2
ATOM 2067 CA SER 297	30.810	10.846	9.234	1.00	46.73	B2	ATOM 2118 HGI THR 303	37.369	17.785	6.178	1.00	0.00	B2
ATOM 2068 CB SER 297	30.239	10.884	7.865	1.00	45.48	B2	ATOM 2119 CG2 THR 303	36.469	19.204	7.927	1.00	38.55	B2
ATOM 2069 CG SER 297	30.988	11.782	7.072	1.00	46.27	B2	ATOM 2120 C THR 303	37.687	18.223	10.505	1.00	34.93	B2
ATOM 2070 HG SER 297	30.321	12.200	6.503	1.00	0.00	B2	ATOM 2121 O THR 303	38.085	19.263	11.063	1.00	35.11	B2
ATOM 2071 C SER 297	32.263	10.269	9.123	1.00	48.72	B2	ATOM 2122 N LEU 304	36.928	17.366	11.150	1.00	33.76	B2
ATOM 2072 O SER 297	33.120	11.122	9.391	1.00	50.55	B2	ATOM 2123 H LEU 304	36.672	16.500	10.762	1.00	0.00	B2
ATOM 2073 N PRO 298	32.655	9.069	8.697	1.00	49.68	B2	ATOM 2124 CA LEU 304	36.436	17.746	12.418	1.00	31.01	B2
ATOM 2074 CD PRO 298	31.782	7.964	8.334	1.00	50.62	B2	ATOM 2125 CB LEU 304	35.345	16.803	12.708	1.00	30.31	B2
ATOM 2075 CA PRO 298	34.049	8.701	8.458	1.00	50.33	B2	ATOM 2126 CG LEU 304	34.234	17.567	13.320	1.00	31.32	B2
ATOM 2076 CB PRO 298	33.948	7.308	7.856	1.00	51.53	B2	ATOM 2127 CD1 LEU 304	33.121	17.626	12.309	1.00	28.87	B2
ATOM 2077 CG PRO 298	32.576	7.266	7.231	1.00	50.43	B2	ATOM 2128 CD2 LEU 304	33.921	16.970	14.692	1.00	34.73	B2
ATOM 2078 C PRO 298	34.795	9.692	7.579	1.00	50.08	B2	ATOM 2129 C LEU 304	37.553	17.726	13.421	1.00	11.86	B2
ATOM 2079 O PRO 298	35.883	10.137	7.930	1.00	50.50	B2	ATOM 2130 O LEU 304	37.615	18.623	14.259	1.00	34.21	B2
ATOM 2080 N GLN 299	34.173	10.086	6.469	1.00	50.48	B2	ATOM 2131 N ASP 305	38.510	16.811	13.326	1.00	30.56	B2
ATOM 2081 H GLN 299	33.279	9.729	6.296	1.00	0.00	B2	ATOM 2132 H ASP 305	39.456	16.117	12.635	1.00	0.00	B2
ATOM 2082 CA GLN 299	34.749	11.050	5.550	1.00	51.74	B2	ATOM 2133 CA ASP 305	39.576	16.797	14.303	1.00	39.72	B2
ATOM 2083 CB GLN 299	33.898	11.236	4.301	1.00	54.33	B2	ATOM 2134 CB ASP 305	40.504	15.608	14.114	1.00	36.20	B2
ATOM 2084 CG GLN 299	33.095	10.067	3.725	1.00	58.11	B2	ATOM 2135 CG ASP 305	39.912	14.201	14.288	1.00	40.64	B2
ATOM 2085 CD GLN 299	31.658	10.086	4.259	1.00	61.49	B2	ATOM 2136 OD1 ASP 305	38.976	14.040	15.103	1.00	37.52	B2
ATOM 2086 OEI GLN 299	31.160	9.083	4.776	1.00	61.00	B2	ATOM 2137 OD2 ASP 305	40.426	13.304	13.581	1.00	42.99	B2
ATOM 2087 NEI GLN 299	30.942	11.217	4.204	1.00	62.12	B2	ATOM 2138 C ASP 305	40.435	18.034	14.238	1.00	27.56	B2
ATOM 2088 HEI2 GLN 299	31.345	12.012	3.800	1.00	0.00	B2	ATOM 2139 O ASP 305	40.775	18.575	15.311	1.00	24.51	B2
ATOM 2089 HEI2 GLN 299	30.034	11.191	4.566	1.00	0.00	B2	ATOM 2140 N THR 306	40.781	18.417	12.979	1.00	24.77	B2

FIGURE 5

ATOM 2141 II THIR 306	40.469	17.875	12.230	1.00	0.00	B2	ATOM 2192 CG1 VAL 311	37.341	25.915	19.949	1.00	19.01	B2
ATOM 2142 CA THIR 306	41.553	19.633	12.751	1.00	24.39	B2	ATOM 2193 CG2 VAL 311	37.261	25.488	17.667	1.00	18.56	B2
ATOM 2143 CB THIR 306	41.465	19.931	11.318	1.00	24.38	B2	ATOM 2194 C VAL 311	40.270	25.638	20.020	1.00	27.21	B2
ATOM 2144 OG1 THIR 306	42.074	18.753	10.665	1.00	25.13	B2	ATOM 2195 O VAL 311	40.437	26.647	20.719	1.00	29.71	B2
ATOM 2145 HG1 THIR 306	41.447	18.029	10.768	1.00	0.00	B2	ATOM 2196 N ALA 312	40.762	24.428	20.357	1.00	27.97	B2
ATOM 2146 CG2 THIR 306	42.690	21.027	11.089	1.00	25.77	B2	ATOM 2197 II ALA 312	40.585	23.674	19.256	1.00	0.00	B2
ATOM 2147 C THIR 306	40.893	20.844	13.419	1.00	25.24	B2	ATOM 2198 CA ALA 312	41.515	24.157	21.583	1.00	27.21	B2
ATOM 2148 O THIR 306	41.488	21.472	14.296	1.00	27.24	B2	ATOM 2199 CB ALA 312	41.855	22.688	21.532	1.00	30.53	B2
ATOM 2149 N LEU 307	39.615	21.134	13.139	1.00	25.91	B2	ATOM 2200 C ALA 312	42.778	25.026	21.784	1.00	30.06	B2
ATOM 2150 II LEU 307	39.125	20.547	12.520	1.00	0.00	B2	ATOM 2201 O ALA 312	43.057	25.508	22.886	1.00	30.04	B2
ATOM 2151 CA LEU 307	38.900	22.218	13.764	1.00	25.53	B2	ATOM 2202 N ASP 313	43.554	25.286	20.735	1.00	31.43	B2
ATOM 2152 CB LEU 307	37.571	22.170	13.142	1.00	25.09	B2	ATOM 2203 II ASP 313	43.433	24.730	19.935	1.00	0.00	B2
ATOM 2153 CG LEU 307	36.530	23.097	13.588	1.00	27.93	B2	ATOM 2204 CA ASP 313	44.610	26.275	20.743	1.00	34.22	B2
ATOM 2154 CD1 LEU 307	37.008	24.515	13.484	1.00	29.87	B2	ATOM 2205 CB ASP 313	45.279	26.512	19.447	1.00	38.87	B2
ATOM 2155 CD2 LEU 307	35.311	22.846	12.728	1.00	28.93	B2	ATOM 2206 CG ASP 313	46.071	25.404	18.866	1.00	44.55	B2
ATOM 2156 C LEU 307	38.850	22.214	15.269	1.00	27.09	B2	ATOM 2207 OD1 ASP 313	46.225	25.439	17.636	1.00	48.67	B2
ATOM 2157 O LEU 307	38.854	23.253	15.925	1.00	30.03	B2	ATOM 2208 OD2 ASP 313	46.531	24.553	19.647	1.00	49.25	B2
ATOM 2158 N GLN 308	38.875	21.044	15.879	1.00	29.09	B2	ATOM 2209 C ASP 313	44.187	27.590	21.059	1.00	35.12	B2
ATOM 2159 H GLN 308	38.883	20.239	15.319	1.00	0.00	B2	ATOM 2210 O ASP 313	44.807	28.390	21.894	1.00	38.60	B2
ATOM 2160 CA GLN 308	38.874	20.848	17.340	1.00	29.36	B2	ATOM 2211 N ASP 314	43.192	28.216	20.339	1.00	33.36	B2
ATOM 2161 CB GLN 308	38.379	19.399	17.562	1.00	29.41	B2	ATOM 2212 II PHE 314	42.784	27.683	19.619	1.00	0.00	B2
ATOM 2162 CG GLN 308	37.862	19.140	18.935	1.00	32.24	B2	ATOM 2213 CA PHE 314	42.715	29.548	20.600	1.00	31.09	B2
ATOM 2163 CD GLN 308	37.586	17.672	19.165	1.00	34.03	B2	ATOM 2214 CB PHE 314	41.572	29.860	19.631	1.00	34.06	B2
ATOM 2164 OE1 GLN 308	36.973	17.023	18.311	1.00	33.48	B2	ATOM 2215 CG PHE 314	41.074	31.303	19.636	1.00	33.37	B2
ATOM 2165 NE2 GLN 308	38.053	17.127	20.299	1.00	31.67	B2	ATOM 2216 CD1 PHE 314	39.780	31.568	19.247	1.00	31.81	B2
ATOM 2166 IE21 GLN 308	38.547	17.697	20.917	1.00	0.00	B2	ATOM 2217 CD2 PHE 314	41.907	32.354	20.021	1.00	35.65	B2
ATOM 2167 IE22 GLN 308	37.875	16.174	20.436	1.00	0.00	B2	ATOM 2218 CE1 PHE 314	39.318	32.857	19.240	1.00	29.15	B2
ATOM 2168 C GLN 308	40.154	21.138	18.051	1.00	28.94	B2	ATOM 2219 CE2 PHE 314	41.455	33.648	20.017	1.00	37.48	B2
ATOM 2169 O GLN 308	40.196	21.796	19.101	1.00	28.44	B2	ATOM 2220 CZ PHE 314	40.154	33.870	19.622	1.00	32.81	B2
ATOM 2170 N LEU 309	41.269	20.671	17.460	1.00	28.78	B2	ATOM 2221 C PHE 314	42.282	29.601	21.057	1.00	29.90	B2
ATOM 2171 II LEU 309	41.157	20.120	16.655	1.00	0.00	B2	ATOM 2222 O PHE 314	42.658	30.550	22.764	1.00	26.87	B2
ATOM 2172 CA LEU 309	42.632	20.923	17.967	1.00	28.56	B2	ATOM 2223 N ALA 315	41.686	28.332	22.584	1.00	29.29	B2
ATOM 2173 CB LEU 309	43.671	20.154	17.106	1.00	26.54	B2	ATOM 2224 H ALA 315	41.448	27.764	22.022	1.00	0.00	B2
ATOM 2174 CG LEU 309	43.632	18.636	17.241	1.00	24.98	B2	ATOM 2225 CA ALA 315	41.300	28.583	23.961	1.00	31.61	B2
ATOM 2175 CD1 LEU 309	44.595	17.935	16.353	1.00	24.17	B2	ATOM 2226 CB ALA 315	40.632	27.358	24.451	1.00	32.23	B2
ATOM 2176 CD2 LEU 309	43.992	18.310	18.621	1.00	23.45	B2	ATOM 2227 C ALA 315	42.482	28.751	24.836	1.00	34.41	B2
ATOM 2177 C LEU 309	42.893	22.416	17.909	1.00	28.24	B2	ATOM 2228 O ALA 315	42.361	29.437	25.853	1.00	37.66	B2
ATOM 2178 O LEU 309	43.370	22.957	18.907	1.00	30.32	B2	ATOM 2229 N THIR 316	43.646	28.250	24.476	1.00	36.16	B2
ATOM 2179 N ASP 310	42.548	23.027	16.749	1.00	26.58	B2	ATOM 2230 II THIR 316	43.745	27.778	23.625	1.00	0.00	B2
ATOM 2180 H ASP 310	42.296	22.437	16.007	1.00	0.00	B2	ATOM 2231 CA THIR 316	44.780	28.388	25.374	1.00	37.99	B2
ATOM 2181 CA ASP 310	42.495	24.477	16.495	1.00	27.90	B2	ATOM 2232 CB THIR 316	45.795	27.255	25.156	1.00	41.16	B2
ATOM 2182 CB ASP 310	42.025	24.659	15.076	1.00	28.41	B2	ATOM 2233 OG1 THIR 316	45.049	26.081	25.521	1.00	45.50	B2
ATOM 2183 CG ASP 310	43.162	24.556	14.096	1.00	31.84	B2	ATOM 2234 HG1 THIR 316	44.316	25.900	24.911	1.00	0.00	B2
ATOM 2184 OD1 ASP 310	42.959	24.766	12.905	1.00	31.54	B2	ATOM 2235 CG2 THIR 316	47.152	27.415	25.888	1.00	40.31	B2
ATOM 2185 OD2 ASP 310	44.297	24.314	14.514	1.00	37.32	B2	ATOM 2236 C THIR 316	45.458	29.710	25.177	1.00	38.47	B2
ATOM 2186 C ASP 310	41.666	25.410	17.422	1.00	27.23	B2	ATOM 2237 O THIR 316	45.903	30.189	26.217	1.00	39.64	B2
ATOM 2187 O ASP 310	42.219	26.429	17.876	1.00	27.23	B2	ATOM 2238 N THIR 317	45.620	30.287	23.970	1.00	36.53	B2
ATOM 2188 N VAL 311	40.374	25.086	17.725	1.00	26.29	B2	ATOM 2239 II THIR 317	45.351	29.800	23.164	1.00	0.00	B2
ATOM 2189 II VAL 311	39.961	24.347	17.725	1.00	0.00	B2	ATOM 2240 CA THIR 317	46.097	31.657	23.844	1.00	37.07	B2
ATOM 2190 CA VAL 311	39.546	25.803	18.706	1.00	24.29	B2	ATOM 2241 CB THIR 317	45.866	32.098	22.392	1.00	36.01	B2
ATOM 2191 CB VAL 311	38.098	25.217	18.869	1.00	21.47	B2	ATOM 2242 OG1 THIR 317	46.752	31.352	21.575	1.00	45.11	B2

FIGURE 5

ATOM 2243	HG1 THR	317	46.409	30.441	21.389	1.00	0.00	B2
ATOM 2244	CG2 THR	317	46.109	33.566	22.156	1.00	34.30	B2
ATOM 2245	C THR	317	45.338	32.597	24.832	1.00	39.30	B2
ATOM 2246	O THR	317	45.941	33.378	25.583	1.00	40.17	B2
ATOM 2247	N ILE	318	44.003	32.481	24.912	1.00	40.83	B2
ATOM 2248	II ILE	318	43.554	31.819	24.342	1.00	0.00	B2
ATOM 2249	CA ILE	318	43.172	33.317	25.788	1.00	40.75	B2
ATOM 2250	CB ILE	318	41.621	32.979	25.567	1.00	37.17	B2
ATOM 2251	CG2 ILE	318	40.742	33.706	26.545	1.00	34.29	B2
ATOM 2252	CG1 ILE	318	41.216	33.310	24.160	1.00	31.39	B2
ATOM 2253	CD ILE	318	41.626	34.657	23.614	1.00	29.66	B2
ATOM 2254	C ILE	318	43.624	33.019	27.217	1.00	42.43	B2
ATOM 2255	O ILE	318	44.064	33.963	27.856	1.00	42.54	B2
ATOM 2256	N TRP	319	43.662	31.784	27.744	1.00	44.17	B2
ATOM 2257	H TRP	319	43.537	31.008	27.163	1.00	0.00	B2
ATOM 2258	CA TRP	319	43.994	31.633	29.142	1.00	46.90	B2
ATOM 2259	CB TRP	319	43.892	30.179	29.597	1.00	50.64	B2
ATOM 2260	CG TRP	319	43.998	30.094	31.131	1.00	56.05	B2
ATOM 2261	CD2 TRP	319	43.005	30.397	32.038	1.00	58.61	B2
ATOM 2262	CE2 TRP	319	43.685	30.281	33.251	1.00	60.50	B2
ATOM 2263	CE3 TRP	319	41.668	30.740	32.005	1.00	60.12	B2
ATOM 2264	CD1 TRP	319	45.188	29.788	31.760	1.00	58.07	B2
ATOM 2265	NE1 TRP	319	44.968	29.921	33.042	1.00	60.07	B2
ATOM 2266	HE1 TRP	319	45.637	29.765	33.740	1.00	0.00	B2
ATOM 2267	CZ2 TRP	319	43.044	30.312	34.456	1.00	61.00	B2
ATOM 2268	CZ3 TRP	319	41.022	30.967	33.210	1.00	61.58	B2
ATOM 2269	CH2 TRP	319	41.704	30.854	34.417	1.00	62.04	B2
ATOM 2270	C TRP	319	45.398	32.136	29.456	1.00	47.85	B2
ATOM 2271	O TRP	319	45.635	32.772	30.490	1.00	47.99	B2
ATOM 2272	N GLN	320	46.339	31.915	28.550	1.00	48.63	B2
ATOM 2273	II GLN	320	46.091	31.482	27.708	1.00	0.00	B2
ATOM 2274	CA GLN	320	47.706	32.319	28.767	1.00	49.45	B2
ATOM 2275	CB GLN	320	48.567	31.988	27.589	1.00	51.44	B2
ATOM 2276	CG GLN	320	48.828	30.494	27.444	1.00	55.03	B2
ATOM 2277	CD GLN	320	49.958	30.349	26.438	1.00	60.17	B2
ATOM 2278	OE1 GLN	320	51.116	30.465	26.834	1.00	65.26	B2
ATOM 2279	NE2 GLN	320	49.771	30.145	25.131	1.00	59.32	B2
ATOM 2280	HE2 GLN	320	48.859	30.087	24.789	1.00	0.00	B2
ATOM 2281	HE22 GLN	320	50.582	30.083	24.590	1.00	0.00	B2
ATOM 2282	C GLN	320	47.717	33.790	28.983	1.00	49.62	B2
ATOM 2283	O GLN	320	48.251	34.209	29.987	1.00	49.91	B2
ATOM 2284	N GLN	321	46.998	34.538	28.150	1.00	51.76	B2
ATOM 2285	H GLN	321	46.335	34.102	27.403	1.00	0.00	B2
ATOM 2286	CA GLN	321	46.837	33.988	28.278	1.00	52.08	B2
ATOM 2287	CB GLN	321	46.015	36.571	27.151	1.00	49.72	B2
ATOM 2288	CG GLN	321	45.873	38.058	27.166	1.00	51.19	B2
ATOM 2289	CD GLN	321	47.211	38.781	27.201	1.00	53.13	B2
ATOM 2290	OE1 GLN	321	48.090	38.622	26.364	1.00	55.36	B2
ATOM 2291	NE2 GLN	321	47.468	39.618	28.177	1.00	53.21	B2
ATOM 2292	HE21 GLN	321	46.800	39.713	28.889	1.00	0.00	B2
ATOM 2293	HE22 GLN	321	48.338	40.057	28.168	1.00	0.00	B2
ATOM 2294	CG PHE	341	46.112	36.315	29.562	1.00	53.30	B2
ATOM 2295	O GLN	321	46.293	37.422	30.058	1.00	54.39	B2
ATOM 2296	N MET	322	45.269	35.441	30.117	1.00	54.50	B2
ATOM 2297	II MET	322	45.098	34.592	29.662	1.00	0.00	B2
ATOM 2298	CA MET	322	44.619	35.748	31.375	1.00	55.42	B2
ATOM 2299	CB MET	322	43.595	34.690	31.713	1.00	52.91	B2
ATOM 2300	CG MET	322	42.527	34.865	30.658	1.00	51.76	B2
ATOM 2301	SD MET	322	40.861	34.428	31.189	1.00	54.19	B2
ATOM 2302	CE MET	322	40.293	33.192	30.069	1.00	52.53	B2
ATOM 2303	C MET	322	45.700	35.811	32.432	1.00	57.69	B2
ATOM 2304	O MET	322	45.781	36.739	33.248	1.00	57.85	B2
ATOM 2305	N GLU	323	46.652	34.900	32.319	1.00	60.28	B2
ATOM 2306	II GLU	323	46.637	34.296	31.544	1.00	0.00	B2
ATOM 2307	CA GLU	323	47.741	34.875	33.273	1.00	62.99	B2
ATOM 2308	CB GLU	323	48.538	33.635	32.957	1.00	65.81	B2
ATOM 2309	CG GLU	323	47.640	32.423	32.918	1.00	68.36	B2
ATOM 2310	CD GLU	323	48.303	31.125	33.310	1.00	71.21	B2
ATOM 2311	OE1 GLU	323	47.651	30.364	34.044	1.00	71.19	B2
ATOM 2312	OE2 GLU	323	49.451	30.900	32.884	1.00	72.43	B2
ATOM 2313	C GLU	323	48.648	36.124	33.418	1.00	63.96	B2
ATOM 2314	OT1 GLU	323	48.782	36.492	34.584	1.00	64.11	B2
ATOM 2315	OT2 GLU	323	49.169	36.725	32.449	1.00	62.96	B2
ATOM 2316	CB MET	338	27.559	17.690	25.056	1.00	62.56	B3
ATOM 2317	CG MET	338	28.087	18.862	24.222	1.00	63.85	B3
ATOM 2318	SD MET	338	28.738	20.224	25.219	1.00	66.95	B3
ATOM 2319	CE MET	338	27.328	21.252	25.515	1.00	65.50	B3
ATOM 2320	C MET	338	24.988	17.301	25.122	1.00	57.55	B3
ATOM 2321	O MET	338	24.417	16.347	25.667	1.00	56.47	B3
ATOM 2322	HT1 MET	338	26.255	16.010	26.594	1.00	0.00	B3
ATOM 2323	HT2 MET	338	25.375	17.061	27.500	1.00	0.00	B3
ATOM 2324	N MET	338	26.286	16.971	27.009	1.00	61.55	B3
ATOM 2325	HT3 MET	338	27.108	17.107	27.620	1.00	0.00	B3
ATOM 2326	CA MET	338	26.226	17.853	25.851	1.00	60.35	B3
ATOM 2327	N PRO	339	24.493	17.830	23.998	1.00	55.58	B3
ATOM 2328	CD PRO	339	24.914	19.075	23.375	1.00	54.39	B3
ATOM 2329	CA PRO	339	23.453	17.226	23.164	1.00	54.62	B3
ATOM 2330	CB PRO	339	23.463	18.098	21.903	1.00	53.52	B3
ATOM 2331	CG PRO	339	24.845	18.711	21.909	1.00	53.04	B3
ATOM 2332	C PRO	339	23.666	15.748	22.881	1.00	53.61	B3
ATOM 2333	O PRO	339	24.730	15.222	23.169	1.00	53.35	B3
ATOM 2334	N ALA	340	22.704	15.045	22.333	1.00	54.32	B3
ATOM 2335	H ALA	340	21.844	15.460	22.111	1.00	0.00	B3
ATOM 2336	CA ALA	340	22.909	13.651	21.968	1.00	56.04	B3
ATOM 2337	CB ALA	340	21.867	12.713	22.625	1.00	57.60	B3
ATOM 2338	C ALA	340	22.617	13.713	20.495	1.00	55.61	B3
ATOM 2339	O ALA	340	21.426	13.783	20.196	1.00	58.64	B3
ATOM 2340	N PHE	341	23.516	13.734	19.514	1.00	53.44	B3
ATOM 2341	II PHE	341	24.472	13.607	19.685	1.00	0.00	B3
ATOM 2342	CA PHE	341	23.016	13.900	18.158	1.00	49.92	B3
ATOM 2343	CB PHE	341	24.050	14.541	17.244	1.00	48.16	B3
ATOM 2344	CG PHE	341	24.382	15.940	17.658	1.00	45.00	B3

FIGURE 5

ATOM 2345 CD1 PHE 341	23.510	16.923	17.359	1.00	43.44	B3
ATOM 2346 CD2 PHE 341	25.527	16.175	18.388	1.00	47.03	B3
ATOM 2347 CE1 PHE 341	23.812	18.172	17.831	1.00	49.15	B3
ATOM 2348 CE2 PHE 341	25.827	17.426	18.862	1.00	47.86	B3
ATOM 2349 CZ PHE 341	24.952	18.437	18.580	1.00	48.36	B3
ATOM 2350 C PHE 341	22.684	12.510	17.672	1.00	49.56	B3
ATOM 2351 O PHE 341	23.309	11.938	16.781	1.00	51.46	B3
ATOM 2352 N ALA 342	21.625	11.985	18.245	1.00	47.40	B3
ATOM 2353 H ALA 342	21.026	12.585	18.741	1.00	47.00	B3
ATOM 2354 CA ALA 342	21.167	10.650	17.997	1.00	46.11	B3
ATOM 2355 CB ALA 342	19.874	10.531	18.804	1.00	47.10	B3
ATOM 2356 C ALA 342	20.962	10.149	16.556	1.00	44.37	B3
ATOM 2357 O ALA 342	20.138	9.747	16.418	1.00	45.65	B3
ATOM 2358 N SER 343	21.537	10.573	15.423	1.00	41.37	B3
ATOM 2359 H SER 343	22.191	11.301	15.428	1.00	41.00	B3
ATOM 2360 CA SER 343	21.274	9.923	14.145	1.00	38.80	B3
ATOM 2361 CB SER 343	19.842	10.138	13.656	1.00	36.79	B3
ATOM 2362 OG SER 343	19.205	11.300	14.182	1.00	37.75	B3
ATOM 2363 HG SER 343	18.963	11.059	15.092	1.00	40.00	B3
ATOM 2364 C SER 343	22.172	10.467	13.088	1.00	38.22	B3
ATOM 2365 O SER 343	22.810	11.471	13.382	1.00	38.30	B3
ATOM 2366 N ALA 344	22.206	9.845	11.888	1.00	36.73	B3
ATOM 2367 H ALA 344	21.762	8.978	11.805	1.00	40.00	B3
ATOM 2368 CA ALA 344	22.914	10.384	10.715	1.00	38.09	B3
ATOM 2369 CB ALA 344	22.583	9.640	9.422	1.00	36.78	B3
ATOM 2370 C ALA 344	22.472	11.842	10.496	1.00	37.72	B3
ATOM 2371 O ALA 344	23.271	12.765	10.676	1.00	36.42	B3
ATOM 2372 N PHE 345	21.194	12.042	10.163	1.00	36.10	B3
ATOM 2373 H PHE 345	20.668	11.298	9.811	1.00	40.00	B3
ATOM 2374 CA PHE 345	20.564	13.338	10.195	1.00	34.69	B3
ATOM 2375 CB PHE 345	19.040	13.254	10.128	1.00	33.24	B3
ATOM 2376 CG PHE 345	18.462	14.636	9.918	1.00	31.72	B3
ATOM 2377 CD1 PHE 345	17.715	15.223	10.505	1.00	26.64	B3
ATOM 2378 CD2 PHE 345	18.767	15.343	8.745	1.00	29.99	B3
ATOM 2379 CE1 PHE 345	17.284	16.503	10.682	1.00	33.56	B3
ATOM 2380 CE2 PHE 345	18.333	16.619	8.537	1.00	30.81	B3
ATOM 2381 CZ PHE 345	17.581	17.201	9.520	1.00	31.44	B3
ATOM 2382 C PHE 345	20.888	14.145	11.458	1.00	35.02	B3
ATOM 2383 O PHE 345	21.246	15.319	11.292	1.00	37.81	B3
ATOM 2384 N GLN 346	20.814	13.688	12.691	1.00	37.53	B3
ATOM 2385 H GLN 346	20.516	12.278	12.894	1.00	40.00	B3
ATOM 2386 CA GLN 346	21.156	14.586	13.758	1.00	33.46	B3
ATOM 2387 CB GLN 346	20.899	13.985	15.061	1.00	33.80	B3
ATOM 2388 CG GLN 346	19.459	14.284	15.174	1.00	35.68	B3
ATOM 2389 CD GLN 346	18.788	13.658	16.344	1.00	38.48	B3
ATOM 2390 OE1 GLN 346	19.358	13.328	17.374	1.00	41.08	B3
ATOM 2391 NE2 GLN 346	17.508	13.463	16.167	1.00	41.08	B3
ATOM 2392 HE1 GLN 346	17.088	13.724	15.323	1.00	40.00	B3
ATOM 2393 HE2 GLN 346	17.026	13.063	16.919	1.00	40.00	B3
ATOM 2394 C GLN 346	22.364	15.051	13.773	1.00	35.73	B3
ATOM 2395 O GLN 346	22.766	16.231	14.051	1.00	38.18	B3
ATOM 2396 N ARG 347	23.507	14.190	13.431	1.00	35.57	B3
ATOM 2397 H ARG 347	23.248	13.289	13.157	1.00	40.00	B3
ATOM 2398 CA ARG 347	24.907	14.538	13.396	1.00	35.95	B3
ATOM 2399 CB ARG 347	25.760	13.236	13.222	1.00	36.20	B3
ATOM 2400 CD ARG 347	26.198	12.549	14.540	1.00	37.41	B3
ATOM 2401 CG ARG 347	26.986	11.246	14.373	1.00	39.70	B3
ATOM 2402 NE ARG 347	26.072	10.167	14.028	1.00	42.18	B3
ATOM 2403 IE ARG 347	25.416	9.893	14.701	1.00	40.00	B3
ATOM 2404 CZ ARG 347	26.071	9.516	12.846	1.00	48.49	B3
ATOM 2405 NH1 ARG 347	26.938	9.802	11.882	1.00	50.22	B3
ATOM 2406 NH11 ARG 347	27.602	10.528	12.031	1.00	40.00	B3
ATOM 2407 NH12 ARG 347	26.905	9.313	11.011	1.00	40.00	B3
ATOM 2408 NH2 ARG 347	25.130	8.608	12.574	1.00	48.46	B3
ATOM 2409 NH21 ARG 347	24.423	8.408	13.252	1.00	40.00	B3
ATOM 2410 NH22 ARG 347	25.126	8.131	11.697	1.00	40.00	B3
ATOM 2411 C ARG 347	25.183	15.544	12.267	1.00	35.54	B3
ATOM 2412 O ARG 347	25.877	16.349	12.445	1.00	36.73	B3
ATOM 2413 N ARG 348	24.611	15.353	11.096	1.00	34.74	B3
ATOM 2414 H ARG 348	24.043	14.559	11.005	1.00	40.00	B3
ATOM 2415 CA ARG 348	24.802	16.225	9.954	1.00	35.24	B3
ATOM 2416 CB ARG 348	24.091	15.623	8.751	1.00	36.76	B3
ATOM 2417 CG ARG 348	24.778	14.303	8.450	1.00	44.22	B3
ATOM 2418 CD ARG 348	24.014	13.379	7.529	1.00	49.23	B3
ATOM 2419 NE ARG 348	24.705	12.090	7.457	1.00	54.27	B3
ATOM 2420 IE ARG 348	25.300	11.836	8.193	1.00	40.00	B3
ATOM 2421 CZ ARG 348	24.557	11.226	6.430	1.00	53.75	B3
ATOM 2422 NH1 ARG 348	23.758	11.479	5.381	1.00	51.85	B3
ATOM 2423 NH11 ARG 348	23.234	12.329	5.339	1.00	40.00	B3
ATOM 2424 NH12 ARG 348	23.680	10.807	4.645	1.00	40.00	B3
ATOM 2425 NH2 ARG 348	25.252	10.083	6.462	1.00	54.51	B3
ATOM 2426 NH21 ARG 348	25.169	9.424	5.714	1.00	40.00	B3
ATOM 2427 NH22 ARG 348	25.860	9.894	7.232	1.00	40.00	B3
ATOM 2428 C ARG 348	24.283	12.629	10.237	1.00	34.80	B3
ATOM 2429 O ARG 348	25.078	18.564	10.219	1.00	35.16	B3
ATOM 2430 N ALA 349	23.008	17.795	10.607	1.00	33.85	B3
ATOM 2431 H ALA 349	22.470	16.984	10.755	1.00	40.00	B3
ATOM 2432 CA ALA 349	22.352	19.083	10.853	1.00	32.96	B3
ATOM 2433 CB ALA 349	20.809	18.894	11.070	1.00	33.46	B3
ATOM 2434 C ALA 349	22.945	19.746	12.083	1.00	31.84	B3
ATOM 2435 O ALA 349	22.981	20.969	12.210	1.00	30.69	B3
ATOM 2436 N GLY 350	23.444	18.954	13.018	1.00	31.30	B3
ATOM 2437 H GLY 350	23.308	17.984	12.976	1.00	40.00	B3
ATOM 2438 CA GLY 350	24.117	19.505	14.181	1.00	31.08	B3
ATOM 2439 C GLY 350	25.462	20.025	13.753	1.00	30.79	B3
ATOM 2440 O GLY 350	25.974	21.010	14.280	1.00	31.38	B3
ATOM 2441 N GLY 351	25.991	19.374	12.731	1.00	30.04	B3
ATOM 2442 H GLY 351	25.546	18.582	12.367	1.00	40.00	B3
ATOM 2443 CA GLY 351	27.263	19.735	12.184	1.00	29.95	B3
ATOM 2444 C GLY 351	27.182	21.097	11.534	1.00	29.25	B3
ATOM 2445 O GLY 351	27.937	21.974	11.919	1.00	28.74	B3
ATOM 2446 N VAL 352	26.336	21.285	10.522	1.00	28.92	B3

FIGURE 5

ATOM 2447 H VAL 352	25.859	20.484	10.214	1.00	0.00	B3
ATOM 2448 CA VAL 352	26.079	22.567	9.881	1.00	28.59	B3
ATOM 2449 CB VAL 352	24.845	22.452	9.004	1.00	28.96	B3
ATOM 2450 CG VAL 352	24.627	23.785	8.346	1.00	30.86	B3
ATOM 2451 CG2 VAL 352	25.021	21.475	7.875	1.00	26.94	B3
ATOM 2452 C VAL 352	25.849	23.709	10.890	1.00	29.79	B3
ATOM 2453 O VAL 352	26.520	24.747	10.853	1.00	31.02	B3
ATOM 2454 N LEU 353	24.923	23.543	11.819	1.00	27.52	B3
ATOM 2455 H LEU 353	24.404	22.709	11.838	1.00	0.00	B3
ATOM 2456 CA LEU 353	24.635	24.548	12.817	1.00	26.18	B3
ATOM 2457 CB LEU 353	23.434	24.113	13.636	1.00	27.87	B3
ATOM 2458 CG LEU 353	22.098	24.034	12.931	1.00	26.54	B3
ATOM 2459 CD1 LEU 353	21.064	23.617	13.924	1.00	25.49	B3
ATOM 2460 CD2 LEU 353	21.750	25.372	12.320	1.00	28.23	B3
ATOM 2461 C LEU 353	25.742	24.905	13.772	1.00	27.17	B3
ATOM 2462 O LEU 353	25.838	26.093	14.088	1.00	28.00	B3
ATOM 2463 N VAL 354	26.339	23.949	14.318	1.00	27.20	B3
ATOM 2464 H VAL 354	26.321	23.006	14.139	1.00	0.00	B3
ATOM 2465 CA VAL 354	27.712	24.212	15.157	1.00	24.62	B3
ATOM 2466 CB VAL 354	28.236	22.910	15.745	1.00	22.01	B3
ATOM 2467 CG1 VAL 354	29.568	22.089	16.406	1.00	19.82	B3
ATOM 2468 CG2 VAL 354	27.276	22.467	16.802	1.00	23.96	B3
ATOM 2469 C VAL 354	28.812	24.893	14.332	1.00	25.46	B3
ATOM 2470 O VAL 354	29.439	25.832	14.798	1.00	26.23	B3
ATOM 2471 N ALA 355	29.059	24.530	13.089	1.00	26.12	B3
ATOM 2472 H ALA 355	28.379	23.745	12.744	1.00	0.00	B3
ATOM 2473 CA ALA 355	30.025	25.180	12.235	1.00	26.34	B3
ATOM 2474 CB ALA 355	30.034	24.591	10.869	1.00	22.08	B3
ATOM 2475 C ALA 355	29.533	26.601	12.096	1.00	28.51	B3
ATOM 2476 O ALA 355	30.315	27.498	12.344	1.00	31.93	B3
ATOM 2477 N SER 356	28.271	26.884	11.802	1.00	30.30	B3
ATOM 2478 H SER 356	27.654	26.134	11.665	1.00	0.00	B3
ATOM 2479 CA SER 356	27.778	28.249	11.625	1.00	31.10	B3
ATOM 2480 CB SER 356	26.401	28.147	11.016	1.00	35.23	B3
ATOM 2481 CG SER 356	25.679	29.380	10.905	1.00	43.82	B3
ATOM 2482 HG SER 356	26.250	30.004	10.429	1.00	0.00	B3
ATOM 2483 C SER 356	27.763	29.095	12.901	1.00	29.75	B3
ATOM 2484 O SER 356	28.115	30.289	12.898	1.00	28.35	B3
ATOM 2485 N HIS 357	27.465	28.464	14.025	1.00	27.82	B3
ATOM 2486 H HIS 357	27.301	27.498	14.019	1.00	0.00	B3
ATOM 2487 CA HIS 357	27.434	29.194	15.259	1.00	26.58	B3
ATOM 2488 CB HIS 357	26.735	28.365	16.305	1.00	25.77	B3
ATOM 2489 CG HIS 357	25.219	28.360	16.063	1.00	27.67	B3
ATOM 2490 CD2 HIS 357	24.563	28.767	14.915	1.00	28.94	B3
ATOM 2491 ND1 HIS 357	24.277	27.963	16.915	1.00	28.43	B3
ATOM 2492 ID1 HIS 357	24.456	27.622	17.828	1.00	0.00	B3
ATOM 2493 CE1 HIS 357	23.112	28.103	16.357	1.00	28.64	B3
ATOM 2494 NE2 HIS 357	23.298	28.589	15.130	1.00	29.48	B3
ATOM 2495 HE2 HIS 357	22.576	28.801	14.495	1.00	0.00	B3
ATOM 2496 C HIS 357	28.852	29.506	15.645	1.00	27.93	B3
ATOM 2497 O HIS 357	29.119	30.606	16.115	1.00	29.15	B3
ATOM 2498 N LEU 358	29.830	28.637	15.383	1.00	28.33	B3
ATOM 2499 H LEU 358	29.624	27.761	14.997	1.00	0.00	B3
ATOM 2500 CA LEU 358	31.211	28.940	15.721	1.00	26.39	B3
ATOM 2501 CB LEU 358	32.030	27.602	15.547	1.00	21.42	B3
ATOM 2502 CG LEU 358	33.457	27.878	15.734	1.00	20.42	B3
ATOM 2503 CD1 LEU 358	33.805	28.078	17.165	1.00	16.79	B3
ATOM 2504 CD2 LEU 358	34.075	26.714	15.072	1.00	22.89	B3
ATOM 2505 C LEU 358	31.721	30.067	14.822	1.00	26.45	B3
ATOM 2506 O LEU 358	31.372	30.939	15.378	1.00	27.12	B3
ATOM 2507 N GIN 359	31.460	30.130	13.518	1.00	26.41	B3
ATOM 2508 H GIN 359	31.048	29.337	13.114	1.00	0.00	B3
ATOM 2509 CA GIN 359	31.863	31.254	12.671	1.00	29.10	B3
ATOM 2510 CB GIN 359	31.204	31.209	11.292	1.00	30.49	B3
ATOM 2511 CG GIN 359	31.395	29.952	10.455	1.00	38.94	B3
ATOM 2512 CD GIN 359	32.842	29.616	10.091	1.00	42.09	B3
ATOM 2513 OE1 GIN 359	33.774	29.979	10.821	1.00	46.15	B3
ATOM 2514 NE2 GIN 359	33.103	28.987	8.967	1.00	41.34	B3
ATOM 2515 HE2 GIN 359	32.341	28.706	8.412	1.00	0.00	B3
ATOM 2516 HE22 GIN 359	34.039	28.838	8.740	1.00	0.00	B3
ATOM 2518 O GIN 359	31.512	32.621	13.254	1.00	29.49	B3
ATOM 2519 N SER 360	32.427	33.427	13.484	1.00	30.06	B3
ATOM 2520 H SER 360	30.201	32.810	13.578	1.00	28.06	B3
ATOM 2521 CA SER 360	29.595	32.061	13.343	1.00	0.00	B3
ATOM 2522 CB SER 360	29.570	34.003	14.071	1.00	27.31	B3
ATOM 2523 CG SER 360	28.121	33.761	14.336	1.00	26.70	B3
ATOM 2524 HG SER 360	27.493	33.539	13.078	1.00	29.72	B3
ATOM 2525 C SER 360	30.202	34.387	15.353	1.00	27.15	B3
ATOM 2526 O SER 360	30.575	35.550	15.498	1.00	26.93	B3
ATOM 2527 N PHE 361	30.383	33.404	16.246	1.00	25.18	B3
ATOM 2528 H PHE 361	30.055	32.499	16.040	1.00	0.00	B3
ATOM 2529 CA PHE 361	31.066	33.626	17.517	1.00	25.20	B3
ATOM 2530 CB PHE 361	31.092	32.335	18.302	1.00	23.40	B3
ATOM 2531 CG PHE 361	31.796	32.394	19.655	1.00	23.63	B3
ATOM 2532 CD1 PHE 361	31.127	32.854	20.777	1.00	22.44	B3
ATOM 2533 CD2 PHE 361	33.098	31.931	19.770	1.00	23.35	B3
ATOM 2534 CE1 PHE 361	31.772	32.834	22.000	1.00	22.78	B3
ATOM 2535 CE2 PHE 361	33.719	31.921	21.002	1.00	21.26	B3
ATOM 2536 CZ PHE 361	33.058	32.308	22.114	1.00	19.54	B3
ATOM 2537 O PHE 361	32.505	34.143	17.385	1.00	26.56	B3
ATOM 2538 N LEU 362	32.914	34.979	18.183	1.00	26.76	B3
ATOM 2539 H LEU 362	33.309	33.645	16.441	1.00	28.17	B3
ATOM 2540 O LEU 362	32.962	32.921	15.874	1.00	0.00	B3
ATOM 2541 CA LEU 362	34.679	34.089	16.222	1.00	28.89	B3
ATOM 2542 CB LEU 362	35.452	34.125	15.338	1.00	28.18	B3
ATOM 2543 CG LEU 362	35.603	31.656	15.781	1.00	29.61	B3
ATOM 2544 CD1 LEU 362	36.306	30.996	14.633	1.00	31.63	B3
ATOM 2545 CD2 LEU 362	36.374	31.453	17.055	1.00	26.38	B3
ATOM 2546 C LEU 362	34.692	35.449	15.536	1.00	29.18	B3
ATOM 2547 O LEU 362	35.649	36.202	15.748	1.00	27.43	B3
ATOM 2548 N GIN 363	33.664	35.763	14.710	1.00	29.54	B3

FIGURE 5

ATOM	2549	II	GLU	363	33.009	35.066	14.495	1.00	0.00	B3
ATOM	2550	CA	GLU	363	33.496	37.090	14.145	1.00	30.30	B3
ATOM	2551	CB	GLU	363	32.357	37.147	13.228	1.00	30.30	B3
ATOM	2552	CG	GLU	363	32.763	36.735	11.849	1.00	38.69	B3
ATOM	2553	CD	GLU	363	33.642	37.662	11.013	1.00	42.62	B3
ATOM	2554	OE1	GLU	363	33.896	37.282	9.860	1.00	46.58	B3
ATOM	2555	OE2	GLU	363	34.051	38.734	11.488	1.00	46.47	B3
ATOM	2556	C	GLU	363	33.229	38.098	15.244	1.00	30.19	B3
ATOM	2557	O	GLU	363	33.837	39.167	15.239	1.00	30.26	B3
ATOM	2558	N	VAL	364	32.397	37.726	16.217	1.00	30.04	B3
ATOM	2559	H	VAL	364	31.888	36.898	16.100	1.00	0.00	B3
ATOM	2560	CA	VAL	364	32.178	38.522	17.400	1.00	31.90	B3
ATOM	2561	CB	VAL	364	31.014	38.021	18.269	1.00	31.41	B3
ATOM	2562	CG1	VAL	364	30.860	38.111	19.562	1.00	30.73	B3
ATOM	2563	CG2	VAL	364	29.750	38.200	17.497	1.00	29.96	B3
ATOM	2564	C	VAL	364	33.402	38.493	18.275	1.00	35.89	B3
ATOM	2565	O	VAL	364	33.683	39.535	18.855	1.00	37.54	B3
ATOM	2566	N	SER	365	34.173	37.421	18.477	1.00	38.25	B3
ATOM	2567	H	SER	365	33.971	36.577	18.030	1.00	0.00	B3
ATOM	2568	CA	SER	365	35.337	37.478	19.375	1.00	39.61	B3
ATOM	2569	CB	SER	365	36.041	36.113	19.555	1.00	43.00	B3
ATOM	2570	CG	SER	365	35.201	34.953	19.575	1.00	46.29	B3
ATOM	2571	HC	SER	365	34.270	35.189	19.644	1.00	0.00	B3
ATOM	2572	C	SER	365	36.398	38.418	18.840	1.00	38.21	B3
ATOM	2573	O	SER	365	37.103	38.989	19.662	1.00	36.91	B3
ATOM	2574	N	TYR	366	36.575	38.540	17.514	1.00	38.00	B3
ATOM	2575	H	TYR	366	36.079	37.945	16.910	1.00	0.00	B3
ATOM	2576	CA	TYR	366	37.568	39.463	16.969	1.00	39.85	B3
ATOM	2577	CB	TYR	366	37.776	39.330	15.436	1.00	38.53	B3
ATOM	2578	CG	TYR	366	38.662	40.447	14.879	1.00	38.21	B3
ATOM	2579	CD1	TYR	366	38.104	41.464	14.129	1.00	37.18	B3
ATOM	2580	CE1	TYR	366	38.918	42.495	13.678	1.00	41.77	B3
ATOM	2581	CD2	TYR	366	40.021	40.443	15.182	1.00	40.21	B3
ATOM	2582	CE2	TYR	366	40.849	41.466	14.739	1.00	40.76	B3
ATOM	2583	CZ	TYR	366	41.151	43.522	13.493	1.00	41.30	B3
ATOM	2584	OH	TYR	366	40.297	42.504	13.976	1.00	42.82	B3
ATOM	2585	H1	TYR	366	40.743	43.977	12.755	1.00	0.00	B3
ATOM	2586	C	TYR	366	37.133	40.893	17.241	1.00	40.55	B3
ATOM	2587	O	TYR	366	37.917	41.647	17.798	1.00	40.92	B3
ATOM	2588	N	ARG	367	35.933	41.309	16.853	1.00	41.88	B3
ATOM	2589	II	ARG	367	35.360	40.682	16.360	1.00	0.00	B3
ATOM	2590	CA	ARG	367	35.442	42.653	17.139	1.00	43.32	B3
ATOM	2591	CB	ARG	367	34.013	42.709	16.650	1.00	46.82	B3
ATOM	2592	CG	ARG	367	33.528	44.130	16.650	1.00	56.74	B3
ATOM	2593	CD	ARG	367	32.089	44.267	16.248	1.00	61.81	B3
ATOM	2594	NE	ARG	367	31.723	45.687	16.229	1.00	66.59	B3
ATOM	2595	IE	ARG	367	32.438	46.356	16.172	1.00	0.00	B3
ATOM	2596	CE1	ARG	367	30.458	46.091	16.308	1.00	69.75	B3
ATOM	2597	NI1	ARG	367	29.448	45.220	16.413	1.00	72.65	B3
ATOM	2598	HI11	ARG	367	29.631	44.236	16.410	1.00	0.00	B3
ATOM	2599	HI12	ARG	367	28.503	45.548	16.445	1.00	0.00	B3
ATOM	2600	NI12	ARG	367	30.160	47.375	16.163	1.00	71.64	B3
ATOM	2601	HI121	ARG	367	29.204	47.865	16.272	1.00	0.00	B3
ATOM	2602	HI122	ARG	367	30.888	48.043	16.013	1.00	0.00	B3
ATOM	2603	C	ARG	367	35.551	43.011	15.615	1.00	40.96	B3
ATOM	2604	O	ARG	367	35.994	44.090	19.012	1.00	41.10	B3
ATOM	2605	N	VAL	368	35.163	42.135	19.542	1.00	39.83	B3
ATOM	2606	II	VAL	368	34.726	41.316	19.217	1.00	0.00	B3
ATOM	2607	CA	VAL	368	35.331	42.497	20.948	1.00	57.33	B3
ATOM	2608	CB	VAL	368	34.748	41.043	21.664	1.00	55.75	B3
ATOM	2609	CG1	VAL	368	35.087	40.867	23.140	1.00	35.10	B3
ATOM	2610	CG2	VAL	368	33.259	41.230	21.586	1.00	33.28	B3
ATOM	2611	C	VAL	368	36.815	42.459	21.223	1.00	48.75	B3
ATOM	2612	O	VAL	368	37.144	43.498	21.772	1.00	40.11	B3
ATOM	2613	N	LEU	369	37.259	41.600	20.845	1.00	49.59	B3
ATOM	2614	II	LEU	369	37.492	40.818	20.308	1.00	0.00	B3
ATOM	2615	CA	LEU	369	39.180	41.780	21.148	1.00	40.05	B3
ATOM	2616	CB	LEU	369	39.984	40.601	20.679	1.00	37.15	B3
ATOM	2617	CG	LEU	369	39.831	39.335	21.426	1.00	37.54	B3
ATOM	2618	CD1	LEU	369	40.349	38.238	20.528	1.00	39.70	B3
ATOM	2619	CD2	LEU	369	40.563	39.394	22.747	1.00	36.86	B3
ATOM	2620	C	LEU	369	39.817	43.031	20.542	1.00	41.98	B3
ATOM	2621	O	LEU	369	40.711	43.654	21.144	1.00	41.30	B3
ATOM	2622	N	ARG	370	39.333	43.413	19.354	1.00	42.80	B3
ATOM	2623	II	ARG	370	38.619	42.884	18.957	1.00	0.00	B3
ATOM	2624	CA	ARG	370	39.819	44.577	18.663	1.00	41.96	B3
ATOM	2625	CB	ARG	370	39.184	44.569	17.316	1.00	42.06	B3
ATOM	2626	CG	ARG	370	39.424	45.719	16.371	1.00	43.93	B3
ATOM	2627	CD	ARG	370	40.894	45.910	16.169	1.00	45.37	B3
ATOM	2628	NE	ARG	370	41.219	46.681	14.976	1.00	48.00	B3
ATOM	2629	IE	ARG	370	40.524	46.867	14.312	1.00	0.00	B3
ATOM	2630	CZ	ARG	370	42.469	47.153	14.791	1.00	48.45	B3
ATOM	2631	NI11	ARG	370	43.443	46.961	15.691	1.00	49.13	B3
ATOM	2632	HI111	ARG	370	43.262	46.456	16.534	1.00	0.00	B3
ATOM	2633	HI112	ARG	370	44.357	47.326	15.520	1.00	0.00	B3
ATOM	2634	NI12	ARG	370	42.821	47.710	13.035	1.00	47.59	B3
ATOM	2635	HI121	ARG	370	42.163	47.785	12.889	1.00	0.00	B3
ATOM	2636	HI122	ARG	370	43.751	48.057	13.516	1.00	0.00	B3
ATOM	2637	C	ARG	370	39.386	45.740	19.558	1.00	49.12	B3
ATOM	2638	O	ARG	370	40.216	46.615	19.826	1.00	49.67	B3
ATOM	2639	N	IIS	371	38.162	45.728	20.123	1.00	52.30	B3
ATOM	2640	II	IIS	371	37.581	44.955	19.949	1.00	0.00	B3
ATOM	2641	CA	IIS	371	37.745	46.738	21.080	1.00	56.65	B3
ATOM	2642	CB	IIS	371	36.284	46.604	21.459	1.00	62.15	B3
ATOM	2643	CG	IIS	371	35.320	46.991	20.346	1.00	71.70	B3
ATOM	2644	CD2	IIS	371	35.596	47.877	19.313	1.00	75.03	B3
ATOM	2645	ND1	IIS	371	34.067	46.346	20.166	1.00	75.91	B3
ATOM	2646	HDI	IIS	371	33.594	45.817	20.732	1.00	0.00	B3
ATOM	2647	CE1	IIS	371	33.580	47.116	19.077	1.00	77.00	B3
ATOM	2648	NE2	IIS	371	34.507	47.914	18.573	1.00	77.52	B3
ATOM	2649	IE2	IIS	371	34.401	48.460	17.764	1.00	0.00	B3
ATOM	2650	C	IIS	371	38.533	46.669	22.887	1.00	56.97	B3

FIGURE 5

ATOM 2651	O	11S	371	38.458	47.592	23.176	1.00	56.12	B3
ATOM 2652	N	LEU	372	39.271	45.632	22.715	1.00	56.98	B3
ATOM 2653	H	LEU	372	39.302	44.855	22.122	1.00	0.00	B3
ATOM 2654	CA	LEU	372	40.048	45.597	23.939	1.00	57.77	B3
ATOM 2655	CB	LEU	372	39.725	44.272	24.633	1.00	52.79	B3
ATOM 2656	CG	LEU	372	38.566	44.144	25.611	1.00	55.87	B3
ATOM 2657	CD1	LEU	372	37.358	44.892	25.123	1.00	55.77	B3
ATOM 2658	CD2	LEU	372	38.211	42.675	25.749	1.00	55.33	B3
ATOM 2659	C	LEU	372	41.554	45.755	23.647	1.00	58.81	B3
ATOM 2660	O	LEU	372	42.447	45.475	24.476	1.00	59.12	B3
ATOM 2661	N	ALA	373	41.942	46.168	22.447	1.00	59.27	B3
ATOM 2662	H	ALA	373	41.271	46.255	21.731	1.00	0.00	B3
ATOM 2663	CA	ALA	373	43.336	46.475	22.147	1.00	60.03	B3
ATOM 2664	CB	ALA	373	43.755	45.485	21.021	1.00	59.87	B3
ATOM 2665	C	ALA	373	43.616	47.895	21.762	1.00	61.22	B3
ATOM 2666	OT1	ALA	373	44.798	48.243	21.697	1.00	62.45	B3
ATOM 2667	OT2	ALA	373	42.682	48.700	21.583	1.00	61.55	B3
ATOM 2668	CB	LEU	410	23.866	49.243	1.118	1.00	53.10	C1
ATOM 2669	CG	LEU	410	23.982	47.812	0.738	1.00	51.85	C1
ATOM 2670	CD1	LEU	410	25.074	47.596	-0.330	1.00	51.64	C1
ATOM 2671	CD2	LEU	410	24.125	47.081	2.058	1.00	49.28	C1
ATOM 2672	C	LEU	410	22.381	51.214	1.635	1.00	52.99	C1
ATOM 2673	O	LEU	410	22.242	52.166	0.845	1.00	53.00	C1
ATOM 2674	HT1	LEU	410	22.721	50.836	-0.665	1.00	0.00	C1
ATOM 2675	HT2	LEU	410	21.194	50.178	-0.557	1.00	0.00	C1
ATOM 2676	N	LEU	410	22.198	49.968	-0.415	1.00	54.31	C1
ATOM 2677	HT3	LEU	410	22.529	49.174	-0.998	1.00	0.00	C1
ATOM 2678	CA	LEU	410	22.478	49.815	1.004	1.00	53.64	C1
ATOM 2679	N	PRO	411	22.450	51.433	2.965	1.00	52.95	C1
ATOM 2680	CD	PRO	411	22.466	50.407	4.022	1.00	52.54	C1
ATOM 2681	CA	PRO	411	22.666	52.766	3.548	1.00	53.25	C1
ATOM 2682	CB	PRO	411	22.688	52.541	5.068	1.00	52.85	C1
ATOM 2683	CG	PRO	411	23.163	51.108	5.203	1.00	52.83	C1
ATOM 2684	C	PRO	411	23.958	53.413	3.023	1.00	53.47	C1
ATOM 2685	O	PRO	411	25.073	52.878	3.167	1.00	54.02	C1
ATOM 2686	N	GLN	412	23.787	54.599	2.411	1.00	52.79	C1
ATOM 2687	H	GLN	412	22.863	54.900	2.294	1.00	0.00	C1
ATOM 2688	CA	GLN	412	24.873	55.413	1.871	1.00	50.44	C1
ATOM 2689	CB	GLN	412	24.387	56.762	1.413	1.00	52.47	C1
ATOM 2690	CG	GLN	412	23.364	57.408	0.437	1.00	56.51	C1
ATOM 2691	CD	GLN	412	25.228	56.954	-1.017	1.00	59.40	C1
ATOM 2692	OE1	GLN	412	25.869	57.506	-1.913	1.00	59.67	C1
ATOM 2693	NE2	GLN	412	24.336	56.022	-1.369	1.00	60.12	C1
ATOM 2694	HE1	GLN	412	23.734	55.616	-0.737	1.00	0.00	C1
ATOM 2695	HE2	GLN	412	24.396	55.748	-2.328	1.00	0.00	C1
ATOM 2696	C	GLN	412	25.930	55.646	2.916	1.00	48.22	C1
ATOM 2697	O	GLN	412	27.089	55.591	2.545	1.00	46.78	C1
ATOM 2698	N	SER	413	25.614	55.842	4.201	1.00	47.90	C1
ATOM 2699	H	SER	413	24.693	55.976	4.492	1.00	0.00	C1
ATOM 2700	CA	SER	413	26.696	55.984	5.144	1.00	48.75	C1
ATOM 2701	CB	SER	413	26.261	56.344	6.548	1.00	50.61	C1
ATOM 2702	OG	SER	413	27.378	56.872	7.301	1.00	53.05	C1
ATOM 2703	11G	SER	413	28.178	56.355	7.145	1.00	0.00	C1
ATOM 2704	C	SER	413	27.480	54.684	5.267	1.00	48.71	C1
ATOM 2705	O	SER	413	28.698	54.839	5.392	1.00	50.77	C1
ATOM 2706	N	PIE	414	26.947	53.440	5.208	1.00	46.01	C1
ATOM 2707	H	PIE	414	25.906	53.324	5.015	1.00	0.00	C1
ATOM 2708	CA	PIE	414	27.787	52.243	5.274	1.00	42.92	C1
ATOM 2709	CB	PIE	414	26.959	50.915	5.232	1.00	40.76	C1
ATOM 2710	CG	PIE	414	27.633	49.627	4.757	1.00	35.06	C1
ATOM 2711	CD1	PIE	414	27.583	49.256	3.425	1.00	34.71	C1
ATOM 2712	CD2	PIE	414	28.262	48.800	5.663	1.00	34.81	C1
ATOM 2713	CE1	PIE	414	28.156	48.056	3.014	1.00	36.16	C1
ATOM 2714	CE2	PIE	414	28.832	47.602	5.247	1.00	33.46	C1
ATOM 2715	CZ	PIE	414	28.781	47.223	3.923	1.00	34.22	C1
ATOM 2716	C	PIE	414	28.667	52.271	4.044	1.00	41.25	C1
ATOM 2717	O	PIE	414	29.831	51.902	4.110	1.00	41.47	C1
ATOM 2718	N	LEU	415	28.122	52.748	2.942	1.00	39.50	C1
ATOM 2719	H	LEU	415	27.188	53.044	2.946	1.00	0.00	C1
ATOM 2720	CA	LEU	415	28.865	52.769	1.721	1.00	39.91	C1
ATOM 2721	CB	LEU	415	27.946	53.205	0.641	1.00	41.98	C1
ATOM 2722	CG	LEU	415	27.903	52.274	-0.526	1.00	44.75	C1
ATOM 2723	CD1	LEU	415	26.430	51.951	-0.780	1.00	42.94	C1
ATOM 2724	CD2	LEU	415	28.793	52.853	-1.648	1.00	45.94	C1
ATOM 2725	C	LEU	415	30.081	53.669	1.755	1.00	40.03	C1
ATOM 2726	O	LEU	415	31.142	53.348	1.183	1.00	40.28	C1
ATOM 2727	N	LEU	416	29.901	54.779	2.487	1.00	37.46	C1
ATOM 2728	H	LEU	416	29.028	54.948	2.899	1.00	0.00	C1
ATOM 2729	CA	LEU	416	30.942	55.756	2.602	1.00	34.05	C1
ATOM 2730	CB	LEU	416	30.794	57.089	2.908	1.00	34.67	C1
ATOM 2731	CG	LEU	416	29.438	57.704	1.851	1.00	35.24	C1
ATOM 2732	CD1	LEU	416	28.770	58.948	2.358	1.00	31.87	C1
ATOM 2733	CD2	LEU	416	30.310	57.948	0.593	1.00	35.50	C1
ATOM 2734	C	LEU	416	31.952	55.238	3.586	1.00	31.97	C1
ATOM 2735	O	LEU	416	33.131	55.427	3.270	1.00	33.32	C1
ATOM 2736	N	ALA	417	31.573	54.619	4.695	1.00	29.05	C1
ATOM 2737	H	ALA	417	30.621	54.616	4.927	1.00	0.00	C1
ATOM 2738	CA	ALA	417	32.524	53.882	5.561	1.00	29.64	C1
ATOM 2739	CB	ALA	417	31.853	54.087	6.680	1.00	25.16	C1
ATOM 2740	C	ALA	417	33.319	52.827	4.777	1.00	36.68	C1
ATOM 2741	O	ALA	417	34.536	52.721	4.877	1.00	31.52	C1
ATOM 2742	N	CYS	418	32.726	52.041	3.905	1.00	32.19	C1
ATOM 2743	H	CYS	418	31.748	52.017	3.850	1.00	0.00	C1
ATOM 2744	CA	CYS	418	33.499	51.119	3.103	1.00	33.67	C1
ATOM 2745	CB	CYS	418	32.657	50.250	2.226	1.00	33.45	C1
ATOM 2746	CG	CYS	418	31.623	49.208	3.246	1.00	37.80	C1
ATOM 2747	C	CYS	418	34.446	51.818	2.170	1.00	34.80	C1
ATOM 2748	O	CYS	418	35.626	51.441	2.373	1.00	36.47	C1
ATOM 2749	N	LEU	419	34.009	52.820	1.377	1.00	35.49	C1
ATOM 2750	H	LEU	419	33.082	53.131	1.460	1.00	0.00	C1
ATOM 2751	CA	LEU	419	34.886	53.446	0.375	1.00	34.14	C1
ATOM 2752	CB	LEU	419	34.062	54.484	-0.411	1.00	17.09	C1

FIGURE 5

ATOM 2753 CG LEU 419	32.866	53.853	-1.244	1.00	39.61	C1
ATOM 2754 CD1 LEU 419	31.866	54.918	-1.609	1.00	39.24	C1
ATOM 2755 CD2 LEU 419	33.349	53.207	-2.553	1.00	40.02	C1
ATOM 2756 C LEU 419	36.107	54.041	1.047	1.00	31.33	C1
ATOM 2757 O LEU 419	37.198	53.973	0.549	1.00	31.60	C1
ATOM 2758 N GLU 420	35.974	54.483	2.273	1.00	31.92	C1
ATOM 2759 H GLU 420	35.068	54.528	2.648	1.00	0.00	C1
ATOM 2760 CA GLU 420	37.078	54.905	3.092	1.00	31.79	C1
ATOM 2761 CB GLU 420	36.477	55.462	4.344	1.00	34.29	C1
ATOM 2762 CG GLU 420	37.430	56.240	5.185	1.00	38.66	C1
ATOM 2763 CD GLU 420	36.932	56.499	6.609	1.00	45.20	C1
ATOM 2764 OE1 GLU 420	37.873	56.849	7.367	1.00	45.67	C1
ATOM 2765 OE2 GLU 420	35.745	56.345	6.954	1.00	44.21	C1
ATOM 2766 C GLU 420	38.043	53.763	3.423	1.00	31.87	C1
ATOM 2767 O GLU 420	39.253	53.949	3.270	1.00	32.82	C1
ATOM 2768 N GLN 421	37.553	52.624	3.954	1.00	30.46	C1
ATOM 2769 H GLN 421	36.583	52.556	4.098	1.00	0.00	C1
ATOM 2770 CA GLN 421	38.366	51.461	4.783	1.00	29.34	C1
ATOM 2771 CB GLN 421	37.545	50.389	4.984	1.00	30.88	C1
ATOM 2772 CG GLN 421	37.308	50.634	6.463	1.00	33.58	C1
ATOM 2773 CD GLN 421	36.320	49.625	7.058	1.00	37.89	C1
ATOM 2774 OE1 GLN 421	35.357	49.236	6.398	1.00	43.18	C1
ATOM 2775 OE2 GLN 421	36.427	49.095	8.275	1.00	37.13	C1
ATOM 2776 NE1 GLN 421	35.695	48.505	8.536	1.00	0.00	C1
ATOM 2777 HE1 GLN 421	37.207	49.330	8.812	1.00	0.00	C1
ATOM 2778 C GLN 421	38.991	50.862	3.026	1.00	27.36	C1
ATOM 2779 O GLN 421	40.152	50.445	3.099	1.00	29.09	C1
ATOM 2780 N VAL 422	38.379	50.845	1.847	1.00	23.57	C1
ATOM 2781 H VAL 422	37.448	51.138	1.803	1.00	0.00	C1
ATOM 2782 CA VAL 422	39.077	50.420	0.651	1.00	23.52	C1
ATOM 2783 CB VAL 422	38.163	50.636	-0.556	1.00	22.62	C1
ATOM 2784 CG1 VAL 422	38.873	50.455	-1.868	1.00	21.56	C1
ATOM 2785 CG2 VAL 422	37.037	49.610	-0.463	1.00	26.79	C1
ATOM 2786 C VAL 422	40.353	51.254	0.514	1.00	26.22	C1
ATOM 2787 O VAL 422	41.458	50.708	0.508	1.00	28.77	C1
ATOM 2788 N ARG 423	40.275	52.599	0.575	1.00	27.49	C1
ATOM 2789 H ARG 423	39.402	53.016	0.735	1.00	0.00	C1
ATOM 2790 CA ARG 423	41.436	53.456	0.346	1.00	25.91	C1
ATOM 2791 CB ARG 423	41.098	54.943	0.312	1.00	24.39	C1
ATOM 2792 CG ARG 423	40.167	55.366	-0.807	1.00	22.81	C1
ATOM 2793 CD ARG 423	40.525	54.798	-2.172	1.00	25.55	C1
ATOM 2794 NE ARG 423	39.707	55.387	-3.216	1.00	25.38	C1
ATOM 2795 HE ARG 423	39.168	56.173	-2.989	1.00	0.00	C1
ATOM 2796 CZ ARG 423	39.629	54.928	-4.466	1.00	27.32	C1
ATOM 2797 NH1 ARG 423	40.264	53.857	-4.949	1.00	26.37	C1
ATOM 2798 NH11 ARG 423	40.884	53.341	-4.365	1.00	0.00	C1
ATOM 2799 NH12 ARG 423	40.150	53.595	-5.907	1.00	0.00	C1
ATOM 2800 NH2 ARG 423	38.960	55.682	-5.325	1.00	30.38	C1
ATOM 2801 NH21 ARG 423	38.539	56.537	-5.023	1.00	0.00	C1
ATOM 2802 NH22 ARG 423	38.865	55.385	-6.275	1.00	0.00	C1
ATOM 2803 C ARG 423	42.429	53.241	1.432	1.00	23.60	C1
ATOM 2804 O ARG 423	43.594	53.147	1.127	1.00	24.37	C1
ATOM 2805 N LYS 424	42.065	53.050	2.608	1.00	24.38	C1
ATOM 2806 H LYS 424	41.109	53.051	2.890	1.00	0.00	C1
ATOM 2807 CA LYS 424	43.043	52.855	3.722	1.00	25.12	C1
ATOM 2808 CB LYS 424	42.352	52.791	5.051	1.00	23.89	C1
ATOM 2809 CG LYS 424	43.312	52.936	6.190	1.00	28.56	C1
ATOM 2810 CD LYS 424	42.579	52.580	7.486	1.00	35.51	C1
ATOM 2811 CE LYS 424	41.338	53.425	7.853	1.00	40.33	C1
ATOM 2812 NZ LYS 424	40.519	52.722	8.834	1.00	42.73	C1
ATOM 2813 HZ1 LYS 424	41.079	52.559	9.695	1.00	0.00	C1
ATOM 2814 HZ2 LYS 424	40.208	51.814	8.435	1.00	0.00	C1
ATOM 2815 HZ3 LYS 424	39.689	53.306	9.065	1.00	0.00	C1
ATOM 2816 C LYS 424	43.761	51.547	3.462	1.00	27.10	C1
ATOM 2817 O LYS 424	44.923	51.425	3.848	1.00	30.64	C1
ATOM 2818 N ILE 425	42.190	50.542	2.794	1.00	26.83	C1
ATOM 2819 H ILE 425	42.260	50.607	2.488	1.00	0.00	C1
ATOM 2820 CA ILE 425	43.949	49.312	2.561	1.00	25.16	C1
ATOM 2821 CB ILE 425	42.965	48.093	2.336	1.00	24.91	C1
ATOM 2822 CG1 ILE 425	43.654	46.786	1.995	1.00	22.01	C1
ATOM 2823 CG1 ILE 425	42.229	47.909	3.633	1.00	25.14	C1
ATOM 2824 CD ILE 425	40.885	47.169	3.432	1.00	25.68	C1
ATOM 2825 C ILE 425	44.824	49.549	1.346	1.00	23.84	C1
ATOM 2826 O ILE 425	45.959	49.069	1.316	1.00	24.57	C1
ATOM 2827 N GLN 426	44.361	50.267	0.323	1.00	23.28	C1
ATOM 2828 H GLN 426	43.451	50.630	0.393	1.00	0.00	C1
ATOM 2829 CA GLN 426	45.164	50.531	-0.871	1.00	24.13	C1
ATOM 2830 CB GLN 426	44.421	51.344	-1.896	1.00	24.04	C1
ATOM 2831 CG GLN 426	43.275	50.539	-2.396	1.00	23.56	C1
ATOM 2832 CD GLN 426	42.446	51.105	-3.311	1.00	23.92	C1
ATOM 2833 OE1 GLN 426	41.704	52.047	-3.345	1.00	25.44	C1
ATOM 2834 NE2 GLN 426	42.337	50.509	-4.672	1.00	22.55	C1
ATOM 2835 HE1 GLN 426	41.755	50.948	-5.323	1.00	0.00	C1
ATOM 2836 HE2 GLN 426	42.850	49.696	-4.831	1.00	0.00	C1
ATOM 2837 C GLN 426	46.404	51.312	-0.488	1.00	26.69	C1
ATOM 2838 O GLN 426	47.486	51.109	-1.046	1.00	29.73	C1
ATOM 2839 N GLY 427	46.300	52.704	0.499	1.00	26.49	C1
ATOM 2840 H GLY 427	45.410	52.414	0.854	1.00	0.00	C1
ATOM 2841 CA GLY 427	47.446	52.894	1.022	1.00	24.25	C1
ATOM 2842 C GLY 427	48.467	51.913	1.589	1.00	23.08	C1
ATOM 2843 O GLY 427	49.597	51.921	1.106	1.00	22.28	C1
ATOM 2844 N ASP 428	48.107	51.073	2.575	1.00	22.75	C1
ATOM 2845 H ASP 428	47.189	51.111	2.918	1.00	0.00	C1
ATOM 2846 CA ASP 428	49.039	50.108	3.131	1.00	23.87	C1
ATOM 2847 CB ASP 428	48.415	49.199	4.117	1.00	26.52	C1
ATOM 2848 CG ASP 428	47.437	49.779	5.097	1.00	28.84	C1
ATOM 2849 OD1 ASP 428	46.420	49.151	5.265	1.00	31.81	C1
ATOM 2850 OD2 ASP 428	47.662	50.804	5.716	1.00	30.77	C1
ATOM 2851 C ASP 428	49.626	49.191	2.063	1.00	24.16	C1
ATOM 2852 O ASP 428	50.812	48.896	2.088	1.00	26.17	C1
ATOM 2853 N GLY 429	48.840	48.822	1.949	1.00	23.01	C1
ATOM 2854 H GLY 429	47.905	49.113	1.071	1.00	0.00	C1

FIGURE 5

ATOM	2855	CA	GLY	429	49,289	47,964	0.029	1.00	25.44	C1	ATOM	2906	CD	LYS	435	58,244	49,748	4.137	1.00	40.31	C1
ATOM	2856	C	GLY	429	50,405	48,649	-0.716	1.00	27.39	C1	ATOM	2907	CE	LYS	435	58,293	50,861	5.213	1.00	45.32	C1
ATOM	2857	O	GLY	429	51,528	48,135	-0.741	1.00	28.31	C1	ATOM	2908	NZ	LYS	435	58,494	50,325	6.575	1.00	47.31	C1
ATOM	2858	N	ALA	430	50,127	49,840	-1.271	1.00	28.26	C1	ATOM	2909	H21	LYS	435	59,388	49,795	6.611	1.00	0.00	C1
ATOM	2859	H	ALA	430	49,216	50,185	-1.172	1.00	0.00	C1	ATOM	2910	H22	LYS	435	57,708	49,689	6.818	1.00	0.00	C1
ATOM	2860	CA	ALA	430	51,094	50,643	-2.015	1.00	26.04	C1	ATOM	2911	H23	LYS	435	58,534	51,109	7.257	1.00	0.00	C1
ATOM	2861	CB	ALA	430	50,490	51,976	-2.407	1.00	27.93	C1	ATOM	2912	C	LYS	435	59,906	48,135	-0.065	1.00	36.10	C1
ATOM	2862	C	ALA	430	52,300	50,927	-1.133	1.00	25.19	C1	ATOM	2913	O	LYS	435	61,139	48,036	-0.012	1.00	37.08	C1
ATOM	2863	O	ALA	430	53,393	51,053	-1.655	1.00	23.43	C1	ATOM	2914	N	LEU	436	59,215	47,168	-0.665	1.00	36.28	C1
ATOM	2864	N	ALA	431	52,171	50,979	0.186	1.00	24.05	C1	ATOM	2915	H	LEU	436	58,235	47,245	-0.651	1.00	0.00	C1
ATOM	2865	H	ALA	431	51,279	50,872	0.579	1.00	0.00	C1	ATOM	2916	CA	LEU	436	59,793	45,994	-1.304	1.00	34.25	C1
ATOM	2866	CA	ALA	431	53,295	51,213	1.035	1.00	26.29	C1	ATOM	2917	CB	LEU	436	58,655	45,076	-1.753	1.00	34.41	C1
ATOM	2867	CB	ALA	431	52,874	51,522	2.458	1.00	24.14	C1	ATOM	2918	CG	LEU	436	57,970	44,327	-0.610	1.00	34.72	C1
ATOM	2868	C	ALA	431	54,139	49,972	1.073	1.00	29.82	C1	ATOM	2919	CD1	LEU	436	56,764	43,538	-1.181	1.00	33.50	C1
ATOM	2869	O	ALA	431	55,360	50,085	0.959	1.00	31.97	C1	ATOM	2920	CD2	LEU	436	58,880	43,375	0.117	1.00	36.39	C1
ATOM	2870	N	LEU	432	53,562	48,777	1.203	1.00	31.87	C1	ATOM	2921	C	LEU	436	60,669	46,383	-2.467	1.00	33.31	C1
ATOM	2871	H	LEU	432	52,583	48,726	1.279	1.00	0.00	C1	ATOM	2922	O	LEU	436	61,756	45,825	-2.647	1.00	33.94	C1
ATOM	2872	CA	LEU	432	54,337	47,540	1.165	1.00	33.92	C1	ATOM	2923	N	CYS	437	60,220	47,374	-3.222	1.00	32.34	C1
ATOM	2873	CB	LEU	432	53,430	46,315	1.301	1.00	37.42	C1	ATOM	2924	H	CYS	437	59,290	47,661	-3.097	1.00	0.00	C1
ATOM	2874	CG	LEU	432	54,063	44,952	1.574	1.00	37.40	C1	ATOM	2925	CA	CYS	437	60,978	47,949	-4.301	1.00	32.01	C1
ATOM	2875	CD1	LEU	432	54,751	44,949	2.950	1.00	38.10	C1	ATOM	2926	C	CYS	437	62,214	48,704	-3.857	1.00	34.70	C1
ATOM	2876	CD2	LEU	432	52,966	43,901	1.492	1.00	36.27	C1	ATOM	2927	O	CYS	437	63,313	48,599	-4.412	1.00	36.26	C1
ATOM	2877	C	LEU	432	55,096	47,404	-0.146	1.00	33.74	C1	ATOM	2928	CB	CYS	437	60,094	48,840	-5.008	1.00	30.97	C1
ATOM	2878	O	LEU	432	56,306	47,179	-0.138	1.00	33.79	C1	ATOM	2929	SG	CYS	437	61,003	49,666	-6.319	1.00	36.22	C1
ATOM	2879	N	GLN	433	54,402	47,564	-1.276	1.00	34.57	C1	ATOM	2930	N	ALA	438	62,016	49,463	-2.785	1.00	36.35	C1
ATOM	2880	H	GLN	433	53,439	47,733	-1.186	1.00	0.00	C1	ATOM	2931	H	ALA	438	61,108	49,547	-2.431	1.00	0.00	C1
ATOM	2881	CA	GLN	433	55,002	47,526	-2.600	1.00	35.83	C1	ATOM	2932	CA	ALA	438	63,060	50,226	-2.170	1.00	35.83	C1
ATOM	2882	CB	GLN	433	53,999	47,892	-3.664	1.00	35.52	C1	ATOM	2933	CB	ALA	438	62,440	51,107	-1.153	1.00	36.38	C1
ATOM	2883	CG	GLN	433	52,996	46,823	-3.832	1.00	39.40	C1	ATOM	2934	C	ALA	438	64,065	49,794	-1.527	1.00	37.01	C1
ATOM	2884	CD	GLN	433	52,049	47,097	-4.973	1.00	42.46	C1	ATOM	2935	O	ALA	438	65,132	49,168	-2.092	1.00	39.39	C1
ATOM	2885	OEL	GLN	433	50,924	47,526	-4.786	1.00	48.22	C1	ATOM	2936	N	THR	439	63,808	48,591	-0.422	1.00	36.59	C1
ATOM	2886	NE2	GLN	433	52,376	46,878	-6.225	1.00	44.77	C1	ATOM	2937	H	THR	439	62,947	48,723	0.014	1.00	0.00	C1
ATOM	2887	HE1	GLN	433	53,271	46,540	-6.433	1.00	0.00	C1	ATOM	2938	CA	THR	439	64,742	47,669	0.223	1.00	35.70	C1
ATOM	2888	HE2	GLN	433	51,693	47,087	-6.892	1.00	0.00	C1	ATOM	2939	CB	THR	439	64,073	47,042	1.400	1.00	35.34	C1
ATOM	2889	C	GLN	433	56,177	48,485	-2.757	1.00	36.48	C1	ATOM	2940	OG1	THR	439	63,323	48,048	2.040	1.00	38.51	C1
ATOM	2890	O	GLN	433	57,214	48,118	-3.312	1.00	38.08	C1	ATOM	2941	HG1	THR	439	62,419	47,999	1.706	1.00	0.00	C1
ATOM	2891	N	GLU	434	56,055	49,719	-2.287	1.00	36.11	C1	ATOM	2942	CG2	THR	439	65,039	46,479	2.469	1.00	36.50	C1
ATOM	2892	H	GLU	434	55,210	49,978	-1.854	1.00	0.00	C1	ATOM	2943	C	THR	439	65,331	46,517	-0.590	1.00	36.10	C1
ATOM	2893	CA	GLU	434	57,089	50,719	-2.426	1.00	35.93	C1	ATOM	2944	O	THR	439	66,448	46,093	-0.312	1.00	36.51	C1
ATOM	2894	CB	GLU	434	56,408	52,030	-2.068	1.00	41.28	C1	ATOM	2945	N	TYR	440	64,603	45,917	-1.548	1.00	36.02	C1
ATOM	2895	CG	GLU	434	57,126	53,356	-2.019	1.00	43.07	C1	ATOM	2946	H	TYR	440	63,751	46,319	-1.822	1.00	0.00	C1
ATOM	2896	CD	GLU	434	57,832	53,516	-0.698	1.00	45.70	C1	ATOM	2947	CA	TYR	440	65,057	44,691	-2.198	1.00	34.74	C1
ATOM	2897	OEL	GLU	434	57,190	53,538	0.367	1.00	49.33	C1	ATOM	2948	CB	TYR	440	64,175	43,480	-1.878	1.00	33.99	C1
ATOM	2898	OEL	GLU	434	59,051	53,579	-0.760	1.00	45.45	C1	ATOM	2949	CG	TYR	440	64,016	43,240	-0.397	1.00	34.14	C1
ATOM	2899	O	GLU	434	58,257	50,348	-1.548	1.00	34.00	C1	ATOM	2950	CD1	TYR	440	62,773	43,240	0.169	1.00	35.16	C1
ATOM	2900	C	GLU	434	59,388	50,481	-1.983	1.00	32.93	C1	ATOM	2951	CE1	TYR	440	62,635	43,017	1.532	1.00	36.44	C1
ATOM	2901	N	LYS	435	58,067	49,860	-0.330	1.00	34.34	C1	ATOM	2952	CD2	TYR	440	65,126	43,084	0.385	1.00	37.83	C1
ATOM	2902	H	LYS	435	57,146	49,837	0.014	1.00	0.00	C1	ATOM	2953	CE2	TYR	440	64,992	42,881	1.752	1.00	39.02	C1
ATOM	2903	CA	LYS	435	59,151	49,358	0.511	1.00	34.56	C1	ATOM	2954	CZ	TYR	440	63,747	42,864	2.317	1.00	37.34	C1
ATOM	2904	CB	LYS	435	58,577	49,010	1.847	1.00	33.89	C1	ATOM	2955	OH	TYR	440	63,637	42,649	3.678	1.00	37.56	C1
ATOM	2905	CI	LYS	435	58,357	50,231	2.709	1.00	36.71	C1	ATOM	2956	HH	TYR	440	64,498	42,343	3.988	1.00	4.00	C1

FIGURE 5

ATOM 2957 C TYR 440	65.088	44.768	-3.681	1.00	34.07	C1
ATOM 2958 O TYR 440	65.598	43.823	-4.267	1.00	35.54	C1
ATOM 2959 N LYS 441	64.627	45.833	-4.330	1.00	33.18	C1
ATOM 2960 CA LYS 441	64.345	46.623	-3.822	1.00	0.00	C1
ATOM 2961 C LYS 441	64.595	45.957	-5.763	1.00	30.44	C1
ATOM 2962 CB LYS 441	65.983	45.759	-6.364	1.00	33.76	C1
ATOM 2963 CG LYS 441	66.729	47.080	-6.407	1.00	39.59	C1
ATOM 2964 CD LYS 441	67.273	47.497	-5.045	1.00	47.69	C1
ATOM 2965 CE LYS 441	67.503	49.028	-4.984	1.00	53.37	C1
ATOM 2966 NZ LYS 441	65.368	49.549	-4.506	1.00	0.00	C1
ATOM 2967 HZ1 LYS 441	66.267	49.780	-5.240	1.00	0.00	C1
ATOM 2968 HZ2 LYS 441	65.885	49.525	-6.173	1.00	0.00	C1
ATOM 2969 HZ3 LYS 441	66.468	50.801	-5.219	1.00	0.00	C1
ATOM 2970 C LYS 441	63.629	45.015	-6.425	1.00	28.86	C1
ATOM 2971 O LYS 441	63.791	44.688	-7.603	1.00	29.95	C1
ATOM 2972 N LEU 442	62.556	44.601	-5.749	1.00	27.58	C1
ATOM 2973 H LEU 442	62.392	44.924	-4.837	1.00	0.00	C1
ATOM 2974 CA LEU 442	61.554	43.780	-6.402	1.00	28.82	C1
ATOM 2975 CB LEU 442	60.947	41.694	-5.466	1.00	26.98	C1
ATOM 2976 CG LEU 442	61.905	41.634	-4.847	1.00	27.75	C1
ATOM 2977 CD1 LEU 442	61.133	40.643	-4.009	1.00	74.29	C1
ATOM 2978 CD2 LEU 442	62.667	40.932	-5.963	1.00	19.72	C1
ATOM 2979 C LEU 442	60.575	44.892	-6.635	1.00	30.59	C1
ATOM 2980 O LEU 442	59.811	45.261	-5.741	1.00	32.36	C1
ATOM 2981 N CYS 443	60.700	45.506	-7.804	1.00	32.15	C1
ATOM 2982 H CYS 443	61.423	45.199	-8.389	1.00	0.00	C1
ATOM 2983 CA CYS 443	59.866	46.645	-8.191	1.00	32.69	C1
ATOM 2984 C CYS 443	58.807	46.380	-9.217	1.00	33.43	C1
ATOM 2985 O CYS 443	58.051	47.288	-9.465	1.00	34.10	C1
ATOM 2986 CB CYS 443	60.715	47.800	-8.743	1.00	30.74	C1
ATOM 2987 SG CYS 443	61.938	48.345	-7.519	1.00	32.96	C1
ATOM 2988 N HIS 444	58.649	45.260	-9.911	1.00	35.65	C1
ATOM 2989 H HIS 444	59.147	44.445	-9.659	1.00	0.00	C1
ATOM 2990 CA HIS 444	57.662	45.172	-10.975	1.00	37.75	C1
ATOM 2991 CB HIS 444	58.329	45.224	-12.330	1.00	37.09	C1
ATOM 2992 CG HIS 444	59.149	46.476	-12.560	1.00	41.36	C1
ATOM 2993 CD1 HIS 444	60.434	46.664	-12.075	1.00	41.40	C1
ATOM 2994 ND1 HIS 444	58.811	47.563	-13.261	1.00	41.74	C1
ATOM 2995 HD1 HIS 444	57.892	47.890	-13.410	1.00	0.00	C1
ATOM 2996 CE1 HIS 444	59.850	48.372	-13.217	1.00	42.00	C1
ATOM 2997 NE2 HIS 444	60.817	47.832	-12.502	1.00	41.38	C1
ATOM 2998 HE2 HIS 444	61.690	48.748	-12.334	1.00	0.00	C1
ATOM 2999 C HIS 444	56.889	43.871	-10.878	1.00	40.10	C1
ATOM 3000 O HIS 444	57.461	42.867	-11.309	1.00	40.15	C1
ATOM 3001 N PRO 445	55.615	43.752	-10.406	1.00	42.06	C1
ATOM 3002 CD PRO 445	54.738	44.836	-9.937	1.00	41.56	C1
ATOM 3003 CA PRO 445	54.913	42.497	-10.276	1.00	40.90	C1
ATOM 3004 CB PRO 445	53.569	42.882	-9.730	1.00	39.35	C1
ATOM 3005 CG PRO 445	53.364	44.274	-10.215	1.00	39.35	C1
ATOM 3006 H PRO 445	54.868	41.782	-11.600	1.00	42.18	C1
ATOM 3007 C LEU 451	54.769	40.571	-11.569	1.00	45.69	C1
ATOM 3008 N LEU 451	55.082	42.380	-12.769	1.00	41.64	C1
ATOM 3009 H LEU 451	55.320	43.320	-12.761	1.00	0.00	C1
ATOM 3010 CA LEU 451	55.025	41.656	-14.029	1.00	42.05	C1
ATOM 3011 CB LEU 451	54.967	42.639	-15.183	1.00	47.06	C1
ATOM 3012 CG LEU 451	54.109	43.925	-14.992	1.00	56.71	C1
ATOM 3013 CD LEU 451	54.728	45.083	-14.162	1.00	62.28	C1
ATOM 3014 OE1 LEU 451	54.100	45.472	-13.178	1.00	66.26	C1
ATOM 3015 OE2 LEU 451	55.818	45.604	-14.473	1.00	65.55	C1
ATOM 3016 C LEU 451	56.237	40.722	-14.197	1.00	40.44	C1
ATOM 3017 O LEU 451	56.186	39.708	-14.904	1.00	41.66	C1
ATOM 3018 N LEU 451	57.360	40.995	-13.538	1.00	37.89	C1
ATOM 3019 H LEU 451	57.394	41.809	-12.999	1.00	0.00	C1
ATOM 3020 CA LEU 451	58.519	40.096	-13.509	1.00	36.73	C1
ATOM 3021 CB LEU 451	59.750	40.810	-12.976	1.00	44.60	C1
ATOM 3022 CG LEU 451	60.320	41.883	-13.850	1.00	35.27	C1
ATOM 3023 CD LEU 451	61.450	42.699	-13.197	1.00	36.14	C1
ATOM 3024 OE1 LEU 451	62.240	43.286	-13.939	1.00	37.31	C1
ATOM 3025 OE2 LEU 451	61.541	42.782	-11.970	1.00	32.80	C1
ATOM 3026 C LEU 451	58.311	38.850	-12.599	1.00	36.31	C1
ATOM 3027 O LEU 451	59.113	37.911	-12.592	1.00	36.33	C1
ATOM 3028 N LEU 451	57.273	38.765	-11.769	1.00	33.81	C1
ATOM 3029 H LEU 451	56.554	39.431	-11.802	1.00	0.00	C1
ATOM 3030 CA LEU 451	57.145	37.691	-10.849	1.00	31.88	C1
ATOM 3031 CB LEU 451	57.080	38.299	-9.484	1.00	29.29	C1
ATOM 3032 CG LEU 451	58.008	39.432	-9.140	1.00	29.81	C1
ATOM 3033 CD1 LEU 451	57.907	39.863	-7.684	1.00	26.02	C1
ATOM 3034 CD2 LEU 451	59.396	38.931	-9.392	1.00	31.13	C1
ATOM 3035 C LEU 451	55.863	36.977	-11.165	1.00	33.75	C1
ATOM 3036 O LEU 451	55.436	36.145	-10.382	1.00	33.96	C1
ATOM 3037 N VAL 449	55.166	37.233	-12.263	1.00	36.99	C1
ATOM 3038 H VAL 449	55.580	37.800	-12.942	1.00	0.00	C1
ATOM 3039 CA VAL 449	53.819	36.701	-12.472	1.00	41.46	C1
ATOM 3040 CB VAL 449	53.157	37.546	-13.625	1.00	41.56	C1
ATOM 3041 CG1 VAL 449	54.002	37.614	-14.880	1.00	42.22	C1
ATOM 3042 CG2 VAL 449	51.921	36.858	-14.112	1.00	42.01	C1
ATOM 3043 C VAL 449	53.760	35.192	-12.733	1.00	44.81	C1
ATOM 3044 O VAL 449	52.866	34.469	-12.227	1.00	44.54	C1
ATOM 3045 N LEU 450	54.716	34.669	-13.515	1.00	47.21	C1
ATOM 3046 H LEU 450	55.416	35.260	-13.870	1.00	0.00	C1
ATOM 3047 CA LEU 450	54.771	33.243	-13.781	1.00	50.57	C1
ATOM 3048 CB LEU 450	55.942	32.894	-14.678	1.00	50.75	C1
ATOM 3049 CG LEU 450	56.148	33.488	-15.994	1.00	52.39	C1
ATOM 3050 CD1 LEU 450	57.152	32.586	-16.673	1.00	53.05	C1
ATOM 3051 CD2 LEU 450	54.882	33.534	-16.833	1.00	54.10	C1
ATOM 3052 C LEU 450	54.911	32.468	-17.471	1.00	53.83	C1
ATOM 3053 O LEU 450	54.297	31.406	-17.266	1.00	55.62	C1
ATOM 3054 N LEU 451	55.685	33.097	-11.575	1.00	55.46	C1
ATOM 3055 H LEU 451	56.073	33.954	-11.849	1.00	0.00	C1
ATOM 3056 CA LEU 451	55.998	32.654	-10.273	1.00	56.01	C1
ATOM 3057 CB LEU 451	57.137	33.542	-9.731	1.00	55.80	C1
ATOM 3058 CG LEU 451	57.745	33.278	-8.394	1.00	56.96	C1

FIGURE 5

ATOM 3059	CD1 LEU	451	58.833	32.284	-8.653	1.00	59.12	CI	ATOM 3110	C ILE	457	47.048	31.608	-5.472	1.00	81.70	CI
ATOM 3060	CD2 LEU	451	58.369	34.511	-7.751	1.00	58.27	CI	ATOM 3111	O ILE	457	46.903	31.761	-6.700	1.00	82.57	CI
ATOM 3061	C LEU	451	54.785	32.700	-9.280	1.00	55.96	CI	ATOM 3112	N PRO	458	45.963	31.583	-4.705	1.00	81.40	CI
ATOM 3062	O LEU	451	54.717	31.935	-8.319	1.00	53.74	CI	ATOM 3113	CD PRO	458	45.959	31.225	-3.278	1.00	81.21	CI
ATOM 3063	N GLY	452	53.774	33.533	-9.522	1.00	57.52	CI	ATOM 3114	CA PRO	458	44.607	31.643	-5.264	1.00	80.74	CI
ATOM 3064	H GLY	452	53.889	34.241	-10.191	1.00	0.00	CI	ATOM 3115	CB PRO	458	43.779	30.942	-4.157	1.00	81.12	CI
ATOM 3065	CA GLY	452	52.567	33.515	-8.710	1.00	60.66	CI	ATOM 3116	CG PRO	458	44.257	30.293	-3.173	1.00	80.47	CI
ATOM 3066	C GLY	452	51.942	32.137	-8.772	1.00	63.64	CI	ATOM 3117	C PRO	458	44.120	33.083	-5.048	1.00	79.10	CI
ATOM 3067	O GLY	452	51.476	31.593	-7.782	1.00	62.60	CI	ATOM 3118	O PRO	458	43.674	33.736	-4.718	1.00	80.10	CI
ATOM 3068	N HIS	453	52.089	31.545	-9.969	1.00	68.46	CI	ATOM 3119	N TRP	459	44.171	33.662	-6.361	1.00	78.19	CI
ATOM 3069	H HIS	453	52.628	32.040	-10.618	1.00	0.00	CI	ATOM 3120	H TRP	459	44.614	33.185	-7.591	1.00	0.00	CI
ATOM 3070	CA HIS	453	51.606	30.205	-10.326	1.00	72.27	CI	ATOM 3121	CA TRP	459	43.543	34.986	-7.092	1.00	77.73	CI
ATOM 3071	CB HIS	453	51.785	29.908	-11.828	1.00	73.84	CI	ATOM 3122	CB TRP	459	43.802	35.428	-8.522	1.00	78.71	CI
ATOM 3072	CG HIS	453	51.421	31.061	-12.777	1.00	77.81	CI	ATOM 3123	CG TRP	459	43.054	36.677	-5.017	1.00	81.27	CI
ATOM 3073	CD1 HIS	453	50.599	32.148	-12.498	1.00	79.29	CI	ATOM 3124	CD2 TRP	459	41.802	36.771	-9.618	1.00	82.89	CI
ATOM 3074	ND1 HIS	453	51.886	31.244	-14.012	1.00	79.84	CI	ATOM 3125	CE2 TRP	459	41.717	38.139	-9.883	1.00	84.21	CI
ATOM 3075	ND1 HIS	453	52.617	30.739	-14.475	1.00	0.00	CI	ATOM 3126	CE3 TRP	459	40.738	35.960	-9.983	1.00	84.28	CI
ATOM 3076	CE1 HIS	453	51.385	32.382	-14.470	1.00	81.11	CI	ATOM 3127	CD1 TRP	459	43.661	37.899	-8.925	1.00	81.89	CI
ATOM 3077	NE2 HIS	453	50.613	32.923	-13.551	1.00	79.85	CI	ATOM 3128	NE1 TRP	459	42.828	38.765	-9.460	1.00	85.23	CI
ATOM 3078	HE2 HIS	453	50.230	33.825	-13.586	1.00	0.00	CI	ATOM 3129	HE1 TRP	459	42.944	39.738	-9.483	1.00	0.00	CI
ATOM 3079	C HIS	453	52.454	29.235	-9.515	1.00	73.43	CI	ATOM 3130	C22 TRP	459	40.615	38.727	-10.494	1.00	84.17	CI
ATOM 3080	O HIS	453	51.875	28.531	-8.692	1.00	73.36	CI	ATOM 3131	C23 TRP	459	39.630	36.538	-10.597	1.00	84.56	CI
ATOM 3081	N SER	454	53.785	29.207	-9.651	1.00	74.64	CI	ATOM 3132	CH2 TRP	459	39.562	37.904	-10.852	1.00	84.83	CI
ATOM 3082	H SER	454	54.214	29.739	-10.351	1.00	0.00	CI	ATOM 3133	C TRP	459	42.009	35.013	-6.827	1.00	77.31	CI
ATOM 3083	CA SER	454	54.639	28.411	-8.765	1.00	77.07	CI	ATOM 3134	O TRP	459	41.202	34.244	-7.376	1.00	76.38	CI
ATOM 3084	CB SER	454	56.123	28.762	-8.980	1.00	77.34	CI	ATOM 3135	N ALA	460	41.557	35.969	-6.020	1.00	76.81	CI
ATOM 3085	CG SER	454	57.095	27.715	-9.124	1.00	75.28	CI	ATOM 3136	H ALA	460	42.187	36.640	-5.689	1.00	0.00	CI
ATOM 3086	HG SER	454	57.149	27.211	-8.306	1.00	0.00	CI	ATOM 3137	CA ALA	460	40.158	36.044	-5.613	1.00	76.44	CI
ATOM 3087	C SER	454	54.332	28.608	-7.262	1.00	78.84	CI	ATOM 3138	CB ALA	460	40.072	36.724	-4.243	1.00	75.51	CI
ATOM 3088	O SER	454	54.270	27.617	-6.535	1.00	80.57	CI	ATOM 3139	C ALA	460	39.237	36.764	-6.538	1.00	76.29	CI
ATOM 3089	N LEU	455	54.070	29.789	-6.693	1.00	79.72	CI	ATOM 3140	O ALA	460	39.449	37.976	-6.833	1.00	76.98	CI
ATOM 3090	H LEU	455	53.956	30.582	-7.250	1.00	0.00	CI	ATOM 3141	N PRO	461	38.217	36.147	-7.187	1.00	76.26	CI
ATOM 3091	CA LEU	455	53.849	29.915	-5.257	1.00	80.43	CI	ATOM 3142	CD PRO	461	38.104	34.684	-7.245	1.00	75.88	CI
ATOM 3092	CB LEU	455	54.085	31.347	-4.838	1.00	80.20	CI	ATOM 3143	CA PRO	461	37.242	36.793	-8.068	1.00	75.46	CI
ATOM 3093	CG LEU	455	55.389	31.981	-5.269	1.00	81.67	CI	ATOM 3144	CB PRO	461	36.605	35.605	-8.755	1.00	75.71	CI
ATOM 3094	CD1 LEU	455	55.254	33.494	-5.419	1.00	81.36	CI	ATOM 3145	CG PRO	461	36.703	34.458	-7.767	1.00	75.60	CI
ATOM 3095	CD2 LEU	455	56.431	31.579	-4.264	1.00	81.26	CI	ATOM 3146	C PRO	461	36.221	37.803	-7.545	1.00	75.72	CI
ATOM 3096	C LEU	455	52.438	29.510	-4.848	1.00	81.56	CI	ATOM 3147	O PRO	461	35.677	37.734	-6.440	1.00	73.66	CI
ATOM 3097	O LEU	455	52.038	28.893	-3.741	1.00	82.22	CI	ATOM 3148	N LEU	462	35.996	38.767	-8.449	1.00	77.19	CI
ATOM 3098	H LEU	456	51.653	28.816	-5.708	1.00	81.89	CI	ATOM 3149	H LEU	462	36.516	38.723	-9.277	1.00	0.00	CI
ATOM 3099	H GLY	456	52.026	28.592	-6.584	1.00	0.00	CI	ATOM 3150	CA LEU	462	35.069	39.891	-8.275	1.00	78.87	CI
ATOM 3100	CA GLY	456	50.269	28.361	-5.467	1.00	82.22	CI	ATOM 3151	CB LEU	462	35.674	40.984	-7.300	1.00	78.12	CI
ATOM 3101	C GLY	456	49.220	29.386	-4.973	1.00	82.56	CI	ATOM 3152	CG LEU	462	34.786	41.959	-6.558	1.00	78.09	CI
ATOM 3102	O GLY	456	48.268	28.989	-4.276	1.00	82.38	CI	ATOM 3153	CD1 LEU	462	34.051	42.987	-7.406	1.00	78.32	CI
ATOM 3103	N ILE	457	49.342	30.697	-5.286	1.00	82.54	CI	ATOM 3154	CD2 LEU	462	33.767	41.092	-5.828	1.00	78.62	CI
ATOM 3104	H ILE	457	50.075	30.942	-5.894	1.00	0.00	CI	ATOM 3155	C LEU	462	34.701	40.565	-9.611	1.00	80.63	CI
ATOM 3105	CA ILE	457	48.435	31.761	-4.824	1.00	81.63	CI	ATOM 3156	OT1 LEU	462	33.507	40.842	-9.808	1.00	81.74	CI
ATOM 3106	CB ILE	457	49.110	33.157	-5.086	1.00	80.69	CI	ATOM 3157	OT2 LEU	462	35.606	40.847	-10.417	1.00	81.59	CI
ATOM 3107	CG ILE	457	48.218	34.305	-4.662	1.00	79.76	CI	ATOM 3158	CB LEU	472	22.074	42.654	-1.426	1.00	62.24	CI
ATOM 3108	CG ILE	457	50.369	33.275	-4.253	1.00	79.82	CI	ATOM 3159	CG LEU	472	22.278	44.145	-1.189	1.00	59.58	CI
ATOM 3109	CD ILE	457	51.506	33.868	-5.081	1.00	77.89	CI	ATOM 3160	CD1 LEU	472	23.496	44.325	-0.328	1.00	59.11	CI

FIGURE 5

ATOM 3161 CD2 LEU 472	22.501	44.883	-2.486	1.00	56.85	C2
ATOM 3162 C LEU 472	23.504	40.625	-1.996	1.00	63.91	C2
ATOM 3163 O LEU 472	23.738	39.874	-2.949	1.00	64.90	C2
ATOM 3164 HT1 LEU 472	21.563	41.441	-3.595	1.00	0.00	C2
ATOM 3165 HT2 LEU 472	23.091	41.291	-4.237	1.00	0.00	C2
ATOM 3166 N LEU 472	22.472	41.930	-3.693	1.00	64.29	C2
ATOM 3167 HT3 LEU 472	22.358	42.849	-4.160	1.00	0.00	C2
ATOM 3168 CA LEU 472	23.092	42.037	-2.386	1.00	63.85	C2
ATOM 3169 N ALA 473	23.652	40.229	-0.733	1.00	63.02	C2
ATOM 3170 H ALA 473	23.533	40.867	0.002	1.00	0.00	C2
ATOM 3171 CA ALA 473	24.023	38.881	-0.353	1.00	62.37	C2
ATOM 3172 CB ALA 473	22.870	37.939	-0.358	1.00	63.65	C2
ATOM 3173 C ALA 473	25.196	38.354	-1.126	1.00	62.01	C2
ATOM 3174 O ALA 473	26.301	38.651	-0.715	1.00	63.36	C2
ATOM 3175 N GLY 474	25.032	37.784	-2.306	1.00	61.43	C2
ATOM 3176 H GLY 474	24.148	37.818	-2.722	1.00	0.00	C2
ATOM 3177 CA GLY 474	26.101	37.137	-3.047	1.00	63.80	C2
ATOM 3178 C GLY 474	27.354	37.950	-3.356	1.00	65.13	C2
ATOM 3179 O GLY 474	28.482	37.417	-3.257	1.00	66.24	C2
ATOM 3180 N CYS 475	27.175	39.237	-3.757	1.00	64.88	C2
ATOM 3181 H CYS 475	26.261	39.550	-3.885	1.00	0.00	C2
ATOM 3182 CA CYS 475	28.308	40.127	-4.068	1.00	61.84	C2
ATOM 3183 CB CYS 475	27.925	41.413	-4.806	1.00	63.74	C2
ATOM 3184 SG CYS 475	29.494	42.075	-5.437	1.00	68.86	C2
ATOM 3185 C CYS 475	28.995	40.567	-2.795	1.00	57.30	C2
ATOM 3186 O CYS 475	30.214	40.449	-2.724	1.00	57.14	C2
ATOM 3187 N LEU 476	28.230	40.983	-1.779	1.00	53.79	C2
ATOM 3188 H LEU 476	27.264	41.024	-1.885	1.00	0.00	C2
ATOM 3189 CA LEU 476	28.797	41.315	-0.493	1.00	50.43	C2
ATOM 3190 CB LEU 476	27.719	41.723	0.523	1.00	45.68	C2
ATOM 3191 CG LEU 476	27.130	43.165	0.497	1.00	42.80	C2
ATOM 3192 CD1 LEU 476	26.670	43.559	1.896	1.00	36.25	C2
ATOM 3193 CD2 LEU 476	28.180	44.180	0.057	1.00	40.22	C2
ATOM 3194 C LEU 476	29.546	40.108	0.042	1.00	50.42	C2
ATOM 3195 O LEU 476	30.614	40.222	0.646	1.00	50.61	C2
ATOM 3196 N SER 477	29.053	38.922	-0.270	1.00	50.62	C2
ATOM 3197 H SER 477	28.196	38.860	-0.729	1.00	0.00	C2
ATOM 3198 CA SER 477	29.721	37.712	0.125	1.00	51.41	C2
ATOM 3199 CB SER 477	28.778	36.524	-0.031	1.00	53.45	C2
ATOM 3200 OG SER 477	27.732	36.616	0.926	1.00	57.65	C2
ATOM 3201 HC SER 477	27.280	37.462	0.828	1.00	0.00	C2
ATOM 3202 C SER 477	30.978	37.525	-0.681	1.00	50.75	C2
ATOM 3203 O SER 477	31.980	37.143	-0.068	1.00	51.41	C2
ATOM 3204 N GLN 478	31.037	37.788	-1.984	1.00	50.21	C2
ATOM 3205 H GLN 478	30.222	38.056	-2.457	1.00	0.00	C2
ATOM 3206 CA GLN 478	32.307	37.697	-2.715	1.00	51.37	C2
ATOM 3207 CB GLN 478	32.064	37.929	-4.166	1.00	53.65	C2
ATOM 3208 CG GLN 478	31.983	36.570	-4.788	1.00	57.32	C2
ATOM 3209 CD GLN 478	31.354	36.649	-6.160	1.00	60.47	C2
ATOM 3210 OE1 GLN 478	31.999	36.504	-7.205	1.00	62.26	C2
ATOM 3211 NE2 GLN 478	30.045	36.878	-6.167	1.00	62.16	C2
ATOM 3212 HE21 GIN 478	29.569	36.972	-5.317	1.00	0.00	C2
ATOM 3213 HE22 GIN 478	29.641	36.928	-7.054	1.00	0.00	C2
ATOM 3214 C GIN 478	33.398	38.670	-2.249	1.00	50.66	C2
ATOM 3215 O GIN 478	34.584	38.314	-2.217	1.00	50.13	C2
ATOM 3216 N LEU 479	33.045	39.909	-1.859	1.00	48.78	C2
ATOM 3217 H LEU 479	32.131	40.223	-2.039	1.00	0.00	C2
ATOM 3218 CA LEU 479	34.015	40.800	-1.215	1.00	45.87	C2
ATOM 3219 CB LEU 479	33.434	42.141	-0.827	1.00	47.63	C2
ATOM 3220 CG LEU 479	32.853	43.083	-1.818	1.00	49.40	C2
ATOM 3221 CD1 LEU 479	32.596	44.393	-1.078	1.00	48.15	C2
ATOM 3222 CD2 LEU 479	31.779	43.258	-3.000	1.00	48.59	C2
ATOM 3223 C LEU 479	34.505	40.146	0.056	1.00	42.13	C2
ATOM 3224 O LEU 479	35.695	39.955	0.262	1.00	40.50	C2
ATOM 3225 N HIS 480	33.609	39.766	0.950	1.00	49.56	C2
ATOM 3226 H HIS 480	32.658	39.935	0.763	1.00	0.00	C2
ATOM 3227 CA HIS 480	33.979	39.108	2.179	1.00	37.81	C2
ATOM 3228 CB HIS 480	32.742	38.714	2.572	1.00	34.29	C2
ATOM 3229 CG HIS 480	33.094	38.241	4.109	1.00	33.82	C2
ATOM 3230 CD2 HIS 480	33.123	36.932	4.709	1.00	33.44	C2
ATOM 3231 ND1 HIS 480	33.450	38.995	5.344	1.00	34.27	C2
ATOM 3232 HD1 HIS 480	33.505	39.976	5.362	1.00	0.00	C2
ATOM 3233 CE1 HIS 480	33.706	38.223	6.365	1.00	34.80	C2
ATOM 3234 NE2 HIS 480	33.504	36.986	5.965	1.00	37.12	C2
ATOM 3235 HE2 HIS 480	33.637	36.202	6.544	1.00	0.00	C2
ATOM 3236 C HIS 480	34.836	37.860	1.961	1.00	49.08	C2
ATOM 3237 O HIS 480	35.716	37.631	2.791	1.00	40.93	C2
ATOM 3238 N SER 481	34.615	37.029	0.935	1.00	39.24	C2
ATOM 3239 H SER 481	33.900	37.241	0.305	1.00	0.00	C2
ATOM 3240 CA SER 481	35.391	35.818	0.683	1.00	38.17	C2
ATOM 3241 CB SER 481	34.813	34.943	-0.420	1.00	40.42	C2
ATOM 3242 CG SER 481	33.454	34.597	-0.137	1.00	47.61	C2
ATOM 3243 HG SER 481	32.898	35.385	-0.162	1.00	0.00	C2
ATOM 3244 C SER 481	36.724	36.272	0.211	1.00	36.12	C2
ATOM 3245 O SER 481	37.692	35.793	0.265	1.00	36.23	C2
ATOM 3246 N GLY 482	36.786	37.706	-0.744	1.00	36.21	C2
ATOM 3247 H GLY 482	35.956	37.498	-1.168	1.00	0.00	C2
ATOM 3248 CA GLY 482	38.028	37.792	-1.266	1.00	36.50	C2
ATOM 3249 C GLY 482	38.958	38.296	-0.151	1.00	36.14	C2
ATOM 3250 O GLY 482	40.142	37.936	-0.055	1.00	36.65	C2
ATOM 3251 N LEU 483	38.381	39.084	0.750	1.00	44.64	C2
ATOM 3252 H LEU 483	37.445	39.326	0.608	1.00	0.00	C2
ATOM 3253 CA LEU 483	39.073	39.593	1.900	1.00	32.07	C2
ATOM 3254 CB LEU 483	38.134	40.442	2.731	1.00	31.17	C2
ATOM 3255 CG LEU 483	37.535	41.687	2.081	1.00	31.11	C2
ATOM 3256 CD1 LEU 483	36.757	42.411	3.156	1.00	30.82	C2
ATOM 3257 CD2 LEU 483	38.599	42.593	1.480	1.00	24.56	C2
ATOM 3258 C LEU 483	39.600	38.461	2.745	1.00	32.91	C2
ATOM 3259 O LEU 483	40.752	38.498	3.199	1.00	31.25	C2
ATOM 3260 N PHE 484	38.767	37.422	2.925	1.00	34.08	C2
ATOM 3261 H PHE 484	37.907	37.408	2.471	1.00	0.00	C2
ATOM 3262 CA PHE 484	39.105	36.798	3.783	1.00	14.60	C2

FIGURE 5

ATOM 3263 CB PHE 484	37.975	35.300	3.925	1.00	37.46	C2
ATOM 3264 CG PHE 484	38.268	34.183	4.897	1.00	40.86	C2
ATOM 3265 CD1 PHE 484	38.219	32.884	4.482	1.00	45.62	C2
ATOM 3266 CD2 PHE 484	38.528	34.445	6.210	1.00	43.62	C2
ATOM 3267 CE1 PHE 484	38.421	31.858	5.395	1.00	47.98	C2
ATOM 3268 CE2 PHE 484	38.731	33.427	7.119	1.00	46.78	C2
ATOM 3269 C PHE 484	38.677	32.119	6.770	1.00	48.06	C2
ATOM 3270 C PHE 484	40.245	35.602	3.113	1.00	31.92	C2
ATOM 3271 O PHE 484	41.162	35.289	3.826	1.00	34.25	C2
ATOM 3272 N LEU 485	40.376	35.413	1.799	1.00	32.75	C2
ATOM 3273 H LEU 485	39.577	35.717	1.250	1.00	0.00	C2
ATOM 3274 CA LEU 485	41.475	34.778	1.163	1.00	33.74	C2
ATOM 3275 CB LEU 485	41.183	34.629	-0.305	1.00	35.35	C2
ATOM 3276 CG LEU 485	42.101	33.962	-1.275	1.00	37.80	C2
ATOM 3277 CD1 LEU 485	41.181	33.404	-2.345	1.00	41.44	C2
ATOM 3278 CD2 LEU 485	43.125	34.903	-1.899	1.00	40.07	C2
ATOM 3279 C LEU 485	42.740	35.585	1.376	1.00	33.95	C2
ATOM 3280 O LEU 485	43.766	35.060	1.850	1.00	33.84	C2
ATOM 3281 N TYR 486	42.609	36.885	1.034	1.00	33.67	C2
ATOM 3282 H TYR 486	41.757	37.186	0.659	1.00	0.00	C2
ATOM 3283 CA TYR 486	43.662	37.862	1.242	1.00	31.33	C2
ATOM 3284 CB TYR 486	43.210	39.290	0.714	1.00	35.33	C2
ATOM 3285 CG TYR 486	43.300	39.325	-0.825	1.00	33.37	C2
ATOM 3286 CD1 TYR 486	42.154	39.405	-1.579	1.00	32.79	C2
ATOM 3287 CE1 TYR 486	42.228	39.190	-2.944	1.00	33.73	C2
ATOM 3288 CD2 TYR 486	44.533	39.153	-1.445	1.00	34.59	C2
ATOM 3289 CE2 TYR 486	44.618	39.033	-2.818	1.00	34.63	C2
ATOM 3290 CZ TYR 486	43.451	39.096	-3.562	1.00	35.58	C2
ATOM 3291 OH TYR 486	43.484	38.880	-4.942	1.00	38.24	C2
ATOM 3292 HH TYR 486	42.614	39.086	-5.306	1.00	0.00	C2
ATOM 3293 C TYR 486	44.068	37.905	2.697	1.00	27.39	C2
ATOM 3294 O TYR 486	45.258	38.007	2.942	1.00	26.06	C2
ATOM 3295 N GLN 487	43.270	37.691	3.708	1.00	26.95	C2
ATOM 3296 H GLN 487	42.315	37.545	3.565	1.00	0.00	C2
ATOM 3297 CA GLN 487	43.835	37.646	5.031	1.00	28.33	C2
ATOM 3298 CB GLN 487	42.690	37.578	6.050	1.00	32.66	C2
ATOM 3299 CG GLN 487	43.092	37.979	7.485	1.00	37.50	C2
ATOM 3300 CD GLN 487	43.966	39.252	7.469	1.00	40.54	C2
ATOM 3301 OE1 GLN 487	43.441	40.346	7.292	1.00	40.45	C2
ATOM 3302 NE2 GLN 487	45.305	39.206	7.549	1.00	38.19	C2
ATOM 3303 HE1 GLN 487	45.755	40.057	7.452	1.00	0.00	C2
ATOM 3304 HE2 GLN 487	45.736	38.340	7.702	1.00	0.00	C2
ATOM 3305 C GLN 487	44.791	36.455	5.207	1.00	28.53	C2
ATOM 3306 O GLN 487	45.774	36.542	5.964	1.00	28.72	C2
ATOM 3307 N GLY 488	44.550	35.363	4.454	1.00	28.32	C2
ATOM 3308 H GLY 488	43.799	35.400	3.824	1.00	0.00	C2
ATOM 3309 CA GLY 488	45.291	34.120	4.557	1.00	26.04	C2
ATOM 3310 C GLY 488	46.860	34.264	4.033	1.00	25.75	C2
ATOM 3311 O GLY 488	47.660	33.946	4.712	1.00	25.86	C2
ATOM 3312 N LEU 489	46.655	34.798	2.818	1.00	25.05	C2
ATOM 3313 H LEU 489	45.798	35.062	2.416	1.00	0.00	C2
ATOM 3314 CA LEU 489	47.911	34.990	2.099	1.00	25.63	C2
ATOM 3315 CB LEU 489	47.708	35.570	0.725	1.00	27.64	C2
ATOM 3316 CG LEU 489	46.761	34.755	-0.189	1.00	30.83	C2
ATOM 3317 CD1 LEU 489	46.373	35.506	-1.471	1.00	30.43	C2
ATOM 3318 CD2 LEU 489	47.472	33.454	-0.502	1.00	32.62	C2
ATOM 3319 C LEU 489	48.783	35.936	2.853	1.00	25.28	C2
ATOM 3320 O LEU 489	49.973	35.705	2.914	1.00	27.37	C2
ATOM 3321 H LEU 490	48.237	36.935	3.514	1.00	25.79	C2
ATOM 3322 N LEU 490	47.267	37.079	3.515	1.00	0.00	C2
ATOM 3323 CA LEU 490	49.072	37.868	4.220	1.00	25.46	C2
ATOM 3324 CB LEU 490	48.274	39.139	4.567	1.00	27.96	C2
ATOM 3325 CG LEU 490	47.823	40.131	3.474	1.00	27.89	C2
ATOM 3326 CD1 LEU 490	46.772	41.019	4.123	1.00	28.63	C2
ATOM 3327 CD2 LEU 490	48.988	40.942	2.899	1.00	28.15	C2
ATOM 3328 C LEU 490	49.619	37.243	5.459	1.00	27.33	C2
ATOM 3329 O LEU 490	50.740	37.528	5.865	1.00	26.73	C2
ATOM 3330 N GLN 491	48.883	36.370	6.111	1.00	25.88	C2
ATOM 3331 H GLN 491	49.430	36.127	5.799	1.00	0.00	C2
ATOM 3332 CA GLN 491	47.984	35.809	7.314	1.00	33.01	C2
ATOM 3333 CB GLN 491	48.305	35.113	8.027	1.00	38.68	C2
ATOM 3334 CG GLN 491	47.856	35.963	9.197	1.00	46.07	C2
ATOM 3335 CD GLN 491	46.348	36.262	9.278	1.00	50.83	C2
ATOM 3336 OE1 GLN 491	45.965	37.436	9.402	1.00	51.92	C2
ATOM 3337 NE2 GLN 491	45.425	35.794	9.278	1.00	51.67	C2
ATOM 3338 HE1 GLN 491	45.723	34.353	9.286	1.00	0.00	C2
ATOM 3339 HE2 GLN 491	44.489	35.560	9.286	1.00	0.00	C2
ATOM 3340 C GLN 491	50.582	34.867	6.986	1.00	33.58	C2
ATOM 3341 O GLN 491	51.582	34.828	7.715	1.00	34.65	C2
ATOM 3342 N ALA 492	50.482	34.191	5.824	1.00	34.15	C2
ATOM 3343 H ALA 492	49.701	34.382	5.264	1.00	0.00	C2
ATOM 3344 CA ALA 492	51.416	33.177	5.321	1.00	33.64	C2
ATOM 3345 CB ALA 492	52.802	33.678	4.959	1.00	34.79	C2
ATOM 3346 C ALA 492	50.818	32.500	4.081	1.00	31.67	C2
ATOM 3347 O ALA 492	53.789	32.943	4.879	1.00	36.03	C2
ATOM 3348 N LEU 493	52.885	34.981	4.728	1.00	35.94	C2
ATOM 3349 H LEU 493	52.060	35.510	4.721	1.00	0.00	C2
ATOM 3350 CA LEU 493	54.139	35.634	4.426	1.00	34.86	C2
ATOM 3351 CB LEU 493	53.898	36.990	3.747	1.00	31.36	C2
ATOM 3352 CG LEU 493	53.127	37.065	2.443	1.00	28.27	C2
ATOM 3353 CD1 LEU 493	52.715	38.495	2.214	1.00	31.74	C2
ATOM 3354 CD2 LEU 493	53.977	36.608	1.285	1.00	28.79	C2
ATOM 3355 C LEU 493	54.879	35.843	5.721	1.00	36.15	C2
ATOM 3356 O LEU 493	55.985	36.374	5.694	1.00	36.70	C2
ATOM 3357 N GLU 494	54.300	35.497	6.855	1.00	38.35	C2
ATOM 3358 H GLU 494	53.395	35.130	6.836	1.00	0.00	C2
ATOM 3359 CA GLU 494	54.910	35.648	8.157	1.00	43.14	C2
ATOM 3360 CB GLU 494	55.621	34.340	8.545	1.00	46.01	C2
ATOM 3361 CG GLU 494	54.711	33.471	9.419	1.00	53.71	C2
ATOM 3362 CD GLU 494	54.195	32.160	8.785	1.00	60.27	C2
ATOM 3363 OE1 GLU 494	53.146	31.653	9.230	1.00	63.52	C2
ATOM 3364 OE2 GLU 494	54.839	31.630	7.862	1.00	62.76	C2

FIGURE 5

ATOM 3365 C GLU 494	55.865	36.825	8.343	1.00	44.32	C2
ATOM 3366 O GLU 494	57.055	36.678	8.610	1.00	46.91	C2
ATOM 3367 N GLY 495	55.358	38.046	8.114	1.00	44.32	C2
ATOM 3368 H GLY 495	54.450	38.112	7.753	1.00	0.00	C2
ATOM 3369 CA GLY 495	56.104	39.272	8.368	1.00	42.36	C2
ATOM 3370 C GLY 495	57.015	39.695	7.238	1.00	42.33	C2
ATOM 3371 O GLY 495	57.397	40.866	7.220	1.00	42.42	C2
ATOM 3372 N ILE 496	57.310	38.802	6.279	1.00	41.04	C2
ATOM 3373 H ILE 496	56.927	37.906	6.374	1.00	0.00	C2
ATOM 3374 CA ILE 496	58.259	38.993	5.192	1.00	41.15	C2
ATOM 3375 CB ILE 496	57.929	40.216	4.253	1.00	38.60	C2
ATOM 3376 CG2 ILE 496	59.077	40.437	3.248	1.00	37.62	C2
ATOM 3377 CG1 ILE 496	56.662	39.964	3.480	1.00	36.39	C2
ATOM 3378 CD ILE 496	56.314	41.071	2.470	1.00	35.27	C2
ATOM 3379 C ILE 496	59.672	39.203	5.749	1.00	42.91	C2
ATOM 3380 O ILE 496	60.541	38.396	5.448	1.00	44.22	C2
ATOM 3381 N SER 497	59.998	40.278	6.533	1.00	44.31	C2
ATOM 3382 H SER 497	59.297	40.852	6.827	1.00	0.00	C2
ATOM 3383 CA SER 497	61.346	40.501	6.992	1.00	44.86	C2
ATOM 3384 CB SER 497	62.204	41.254	5.938	1.00	40.13	C2
ATOM 3385 CG SER 497	62.181	42.673	6.033	1.00	40.74	C2
ATOM 3386 HG SER 497	62.531	42.964	5.170	1.00	0.00	C2
ATOM 3387 C SER 497	61.164	41.413	8.185	1.00	45.85	C2
ATOM 3388 O SER 497	60.132	42.110	8.288	1.00	47.55	C2
ATOM 3389 N PRO 498	62.164	41.490	9.071	1.00	44.96	C2
ATOM 3390 CD PRO 498	63.338	40.621	9.126	1.00	42.33	C2
ATOM 3391 CA PRO 498	62.086	42.327	10.250	1.00	44.88	C2
ATOM 3392 CB PRO 498	63.431	42.038	10.885	1.00	45.13	C2
ATOM 3393 CG PRO 498	63.629	40.581	10.603	1.00	42.00	C2
ATOM 3394 C PRO 498	61.760	43.799	9.983	1.00	45.22	C2
ATOM 3395 O PRO 498	61.215	44.446	10.869	1.00	45.24	C2
ATOM 3396 N GLU 499	62.017	44.314	8.777	1.00	46.16	C2
ATOM 3397 H GLU 499	62.362	43.716	8.081	1.00	0.00	C2
ATOM 3398 CA GLU 499	61.731	45.699	8.391	1.00	48.06	C2
ATOM 3399 CB GLU 499	62.498	46.193	7.155	1.00	52.19	C2
ATOM 3400 CG GLU 499	64.001	46.187	7.100	1.00	57.51	C2
ATOM 3401 CD GLU 499	64.544	44.727	7.076	1.00	60.61	C2
ATOM 3402 OE1 GLU 499	64.755	44.231	8.162	1.00	62.96	C2
ATOM 3403 OE2 GLU 499	64.739	44.234	5.984	1.00	62.79	C2
ATOM 3404 C GLU 499	60.269	45.896	7.981	1.00	46.94	C2
ATOM 3405 O GLU 499	59.600	46.895	8.272	1.00	48.15	C2
ATOM 3406 N LEU 500	59.806	44.934	7.193	1.00	44.38	C2
ATOM 3407 H LEU 500	60.351	44.137	7.027	1.00	0.00	C2
ATOM 3408 CA LEU 500	58.491	44.997	6.651	1.00	41.08	C2
ATOM 3409 CB LEU 500	58.519	44.197	5.445	1.00	41.37	C2
ATOM 3410 CG LEU 500	59.303	44.862	4.351	1.00	42.70	C2
ATOM 3411 CD1 LEU 500	59.276	43.828	3.351	1.00	43.98	C2
ATOM 3412 CD2 LEU 500	58.427	45.874	3.671	1.00	45.04	C2
ATOM 3413 C LEU 500	57.455	44.521	7.628	1.00	40.59	C2
ATOM 3414 O LEU 500	56.274	44.835	7.463	1.00	40.69	C2
ATOM 3415 N GLY 501	57.866	43.835	8.085	1.00	39.37	C2
ATOM 3416 H GLY 501	58.808	43.579	8.730	1.00	0.00	C2
ATOM 3417 CA GLY 501	56.974	43.386	9.734	1.00	39.59	C2
ATOM 3418 C GLY 501	55.816	44.324	10.092	1.00	39.66	C2
ATOM 3419 O GLY 501	54.661	44.034	9.777	1.00	40.66	C2
ATOM 3420 N PRO 502	55.986	45.462	10.742	1.00	39.90	C2
ATOM 3421 CD PRO 502	57.727	45.908	11.335	1.00	41.18	C2
ATOM 3422 CA PRO 502	54.912	46.387	11.045	1.00	38.67	C2
ATOM 3423 CB PRO 502	55.594	47.494	11.791	1.00	39.23	C2
ATOM 3424 CG PRO 502	56.989	47.405	11.721	1.00	41.36	C2
ATOM 3425 C PRO 502	54.158	46.849	9.817	1.00	37.54	C2
ATOM 3426 O PRO 502	52.966	47.139	9.961	1.00	38.46	C2
ATOM 3427 N THR 503	54.728	46.887	8.609	1.00	35.13	C2
ATOM 3428 H THR 503	55.663	46.638	8.449	1.00	0.00	C2
ATOM 3429 CA THR 503	53.940	47.283	7.462	1.00	35.09	C2
ATOM 3430 CB THR 503	54.832	47.376	6.245	1.00	34.48	C2
ATOM 3431 CG1 THR 503	56.025	48.018	6.668	1.00	38.23	C2
ATOM 3432 HG1 THR 503	55.857	48.946	6.845	1.00	0.00	C2
ATOM 3433 CG2 THR 503	54.197	48.162	5.126	1.00	35.56	C2
ATOM 3434 C THR 503	52.836	46.252	7.215	1.00	35.37	C2
ATOM 3435 O THR 503	51.671	46.552	6.915	1.00	37.11	C2
ATOM 3436 N LEU 504	53.218	44.996	7.380	1.00	34.02	C2
ATOM 3437 H LEU 504	54.146	44.799	7.647	1.00	0.00	C2
ATOM 3438 CA LEU 504	52.301	43.912	7.173	1.00	32.50	C2
ATOM 3439 CB LEU 504	53.127	42.650	7.002	1.00	34.78	C2
ATOM 3440 CG LEU 504	53.464	42.256	5.601	1.00	34.07	C2
ATOM 3441 CD1 LEU 504	54.163	40.977	5.667	1.00	37.97	C2
ATOM 3442 CD2 LEU 504	52.254	41.865	4.809	1.00	37.34	C2
ATOM 3443 C LEU 504	51.324	43.821	8.328	1.00	29.54	C2
ATOM 3444 O LEU 504	50.141	43.562	8.078	1.00	30.40	C2
ATOM 3445 N ASP 505	51.736	44.106	9.551	1.00	26.09	C2
ATOM 3446 H ASP 505	52.689	44.269	9.699	1.00	0.00	C2
ATOM 3447 CA ASP 505	50.798	44.084	10.643	1.00	27.88	C2
ATOM 3448 CB ASP 505	51.446	44.345	11.926	1.00	29.86	C2
ATOM 3449 CG ASP 505	52.500	43.312	12.239	1.00	34.64	C2
ATOM 3450 OD1 ASP 505	52.663	42.298	11.534	1.00	41.04	C2
ATOM 3451 OD2 ASP 505	53.179	43.542	13.224	1.00	37.40	C2
ATOM 3452 C ASP 505	49.661	45.060	10.568	1.00	28.61	C2
ATOM 3453 O ASP 505	48.566	44.739	11.039	1.00	30.30	C2
ATOM 3454 N THR 506	49.894	46.242	10.002	1.00	28.29	C2
ATOM 3455 H THR 506	50.823	46.493	9.804	1.00	0.00	C2
ATOM 3456 CA THR 506	48.860	47.225	9.731	1.00	25.74	C2
ATOM 3457 CB THR 506	49.497	48.556	9.336	1.00	26.14	C2
ATOM 3458 CG1 THR 506	49.944	49.039	10.588	1.00	31.61	C2
ATOM 3459 HG1 THR 506	49.243	49.072	11.246	1.00	0.00	C2
ATOM 3460 CG2 THR 506	48.954	49.517	8.619	1.00	24.46	C2
ATOM 3461 C THR 506	48.022	46.735	8.615	1.00	24.00	C2
ATOM 3462 O THR 506	46.817	46.864	8.719	1.00	23.85	C2
ATOM 3463 N LEU 507	48.554	46.196	7.525	1.00	23.51	C2
ATOM 3464 H LEU 507	49.527	46.073	7.453	1.00	0.00	C2
ATOM 3465 CA LEU 507	47.682	45.770	6.434	1.00	23.85	C2
ATOM 3466 CB LEU 507	48.574	45.408	5.196	1.00	23.33	C2

FIGURE 5

ATOM 3467 CG LEU 507	-8.010	44.919	3.858	1.00	20.85	C2
ATOM 3468 CD1 LEU 507	46.771	45.650	3.455	1.00	24.13	C2
ATOM 3469 CD2 LEU 507	49.074	45.055	2.842	1.00	20.13	C2
ATOM 3470 C LEU 507	46.766	44.640	6.880	1.00	24.09	C2
ATOM 3471 O LEU 507	45.600	44.764	6.541	1.00	25.80	C2
ATOM 3472 N GLN 508	47.152	43.618	7.661	1.00	24.01	C2
ATOM 3473 H GLN 508	48.112	43.555	7.866	1.00	0.00	C2
ATOM 3474 CA GLN 508	46.228	42.625	8.214	1.00	23.71	C2
ATOM 3475 CB GLN 508	46.961	41.627	9.036	1.00	23.83	C2
ATOM 3476 CG GLN 508	47.937	40.899	8.173	1.00	31.64	C2
ATOM 3477 CD GLN 508	48.842	40.080	9.054	1.00	34.00	C2
ATOM 3478 OE1 GLN 508	50.031	40.346	9.161	1.00	38.32	C2
ATOM 3479 NE2 GLN 508	48.321	39.090	9.748	1.00	36.30	C2
ATOM 3480 HE21 GLN 508	47.373	38.880	9.639	1.00	0.00	C2
ATOM 3481 HE22 GLN 508	48.891	38.636	10.406	1.00	0.00	C2
ATOM 3482 C GLN 508	45.105	43.123	9.111	1.00	24.24	C2
ATOM 3483 O GLN 508	43.978	42.650	9.014	1.00	24.06	C2
ATOM 3484 N LEU 509	45.375	44.019	10.090	1.00	26.07	C2
ATOM 3485 H LEU 509	46.316	44.262	10.222	1.00	0.00	C2
ATOM 3486 CA LEU 509	44.378	44.640	10.977	1.00	25.71	C2
ATOM 3487 CB LEU 509	44.993	45.555	12.031	1.00	25.60	C2
ATOM 3488 CG LEU 509	45.838	44.757	13.042	1.00	29.00	C2
ATOM 3489 CD1 LEU 509	46.658	45.705	13.886	1.00	28.93	C2
ATOM 3490 CD2 LEU 509	44.950	43.919	13.937	1.00	27.94	C2
ATOM 3491 C LEU 509	43.465	45.471	10.130	1.00	25.17	C2
ATOM 3492 O LEU 509	42.274	45.411	10.408	1.00	27.22	C2
ATOM 3493 N ASP 510	43.899	46.208	9.101	1.00	23.77	C2
ATOM 3494 H ASP 510	44.865	46.277	8.930	1.00	0.00	C2
ATOM 3495 CA ASP 510	42.955	46.898	8.240	1.00	22.66	C2
ATOM 3496 CB ASP 510	43.652	47.829	7.306	1.00	25.21	C2
ATOM 3497 CG ASP 510	44.316	48.966	8.068	1.00	33.01	C2
ATOM 3498 OD1 ASP 510	45.178	49.621	7.477	1.00	34.28	C2
ATOM 3499 OD2 ASP 510	43.988	49.209	9.250	1.00	34.44	C2
ATOM 3500 C ASP 510	42.104	45.980	7.398	1.00	23.72	C2
ATOM 3501 O ASP 510	40.897	46.220	7.387	1.00	24.80	C2
ATOM 3502 N VAL 511	42.632	44.984	6.659	1.00	22.38	C2
ATOM 3503 H VAL 511	43.611	44.900	6.620	1.00	0.00	C2
ATOM 3504 CA VAL 511	41.823	44.010	5.961	1.00	21.89	C2
ATOM 3505 CB VAL 511	42.752	42.924	5.366	1.00	22.71	C2
ATOM 3506 CG1 VAL 511	41.954	41.756	4.792	1.00	20.43	C2
ATOM 3507 CG2 VAL 511	43.529	43.524	4.210	1.00	16.19	C2
ATOM 3508 C VAL 511	40.827	43.403	6.960	1.00	21.92	C2
ATOM 3509 O VAL 511	39.625	43.447	6.719	1.00	23.46	C2
ATOM 3510 N ALA 512	41.258	43.017	8.163	1.00	20.49	C2
ATOM 3511 H ALA 512	42.216	43.063	8.361	1.00	0.00	C2
ATOM 3512 CA ALA 512	40.388	42.357	9.108	1.00	20.83	C2
ATOM 3513 CB ALA 512	41.103	41.974	10.344	1.00	17.89	C2
ATOM 3514 C ALA 512	39.250	43.205	9.550	1.00	23.89	C2
ATOM 3515 O ALA 512	38.201	42.668	9.874	1.00	24.61	C2
ATOM 3516 N ASP 513	39.417	44.539	9.544	1.00	25.96	C2
ATOM 3517 H ASP 513	40.300	44.888	9.291	1.00	0.00	C2
ATOM 3518 CA ASP 513	38.374	45.471	9.947	1.00	25.37	C2
ATOM 3519 CB ASP 513	38.958	46.787	10.373	1.00	26.88	C2
ATOM 3520 CG ASP 513	39.682	46.679	11.712	1.00	32.35	C2
ATOM 3521 OD1 ASP 513	40.371	47.644	12.058	1.00	35.06	C2
ATOM 3522 OD2 ASP 513	39.580	45.646	12.390	1.00	34.10	C2
ATOM 3523 C ASP 513	37.392	45.730	8.846	1.00	24.95	C2
ATOM 3524 O ASP 513	36.185	45.868	9.090	1.00	26.92	C2
ATOM 3525 N PHE 514	37.867	45.739	7.634	1.00	22.88	C2
ATOM 3526 H PHE 514	38.829	45.614	7.475	1.00	0.00	C2
ATOM 3527 CA PHE 514	36.974	45.922	6.510	1.00	24.09	C2
ATOM 3528 CB PHE 514	37.812	46.061	5.266	1.00	18.11	C2
ATOM 3529 CG PHE 514	36.956	46.470	4.072	1.00	17.86	C2
ATOM 3530 CD1 PHE 514	35.715	47.089	4.245	1.00	15.18	C2
ATOM 3531 CD2 PHE 514	37.440	46.197	2.804	1.00	13.77	C2
ATOM 3532 CE1 PHE 514	34.983	47.419	3.130	1.00	14.53	C2
ATOM 3533 CE2 PHE 514	36.693	46.539	1.705	1.00	12.10	C2
ATOM 3534 C PHE 514	35.468	47.146	1.868	1.00	10.68	C2
ATOM 3535 C PHE 514	36.076	44.703	6.450	1.00	29.23	C2
ATOM 3536 O PHE 514	34.788	44.828	6.350	1.00	29.80	C2
ATOM 3537 N ALA 515	36.604	43.490	6.531	1.00	31.15	C2
ATOM 3538 H ALA 515	37.581	43.450	6.639	1.00	0.00	C2
ATOM 3539 CA ALA 515	35.839	42.260	6.416	1.00	32.36	C2
ATOM 3540 CB ALA 515	36.851	41.126	6.402	1.00	32.35	C2
ATOM 3541 C ALA 515	34.801	42.089	7.535	1.00	32.39	C2
ATOM 3542 O ALA 515	33.676	41.609	7.331	1.00	32.63	C2
ATOM 3543 N THR 516	35.164	42.457	8.235	1.00	33.01	C2
ATOM 3544 H THR 516	36.117	42.578	8.935	1.00	0.00	C2
ATOM 3545 CA THR 516	34.231	42.566	9.821	1.00	35.18	C2
ATOM 3546 CB THR 516	35.016	43.018	10.988	1.00	35.40	C2
ATOM 3547 CG1 THR 516	35.685	41.818	11.336	1.00	42.65	C2
ATOM 3548 HG1 THR 516	36.505	41.713	10.816	1.00	0.00	C2
ATOM 3549 CG2 THR 516	34.262	43.672	12.097	1.00	45.56	C2
ATOM 3550 C THR 516	33.140	43.554	9.482	1.00	37.62	C2
ATOM 3551 O THR 516	32.005	43.315	9.857	1.00	40.37	C2
ATOM 3552 N THR 517	33.387	44.666	8.802	1.00	38.61	C2
ATOM 3553 H THR 517	34.291	44.850	8.469	1.00	0.00	C2
ATOM 3554 CA THR 517	32.359	45.641	8.512	1.00	38.92	C2
ATOM 3555 CB THR 517	33.123	46.903	7.962	1.00	40.46	C2
ATOM 3556 CG1 THR 517	33.832	47.429	9.103	1.00	43.22	C2
ATOM 3557 HG1 THR 517	34.536	46.815	9.335	1.00	0.00	C2
ATOM 3558 CG2 THR 517	32.232	47.926	7.253	1.00	39.90	C2
ATOM 3559 C THR 517	31.343	45.012	7.551	1.00	38.30	C2
ATOM 3560 O THR 517	30.137	45.125	7.811	1.00	38.69	C2
ATOM 3561 N ILE 518	31.790	44.344	6.466	1.00	37.54	C2
ATOM 3562 H ILE 518	32.756	44.386	6.297	1.00	0.00	C2
ATOM 3563 CA ILE 518	30.923	43.646	5.510	1.00	16.10	C2
ATOM 3564 CB ILE 518	31.699	42.912	4.439	1.00	31.81	C2
ATOM 3565 CG1 ILE 518	30.703	42.202	3.555	1.00	13.46	C2
ATOM 3566 CG2 ILE 518	32.623	43.842	3.699	1.00	32.91	C2
ATOM 3567 CD1 ILE 518	32.019	44.700	2.596	1.00	34.89	C2
ATOM 3568 C ILE 518	30.172	42.591	6.317	1.00	38.61	C2

FIGURE 5

ATOM 3569 O HLE 518	28.938	42.545	6.205	1.00	39.93	C2
ATOM 3570 N TRP 519	30.842	41.785	7.179	1.00	38.64	C2
ATOM 3571 H TRP 519	31.785	41.959	7.361	1.00	0.00	C2
ATOM 3572 CA TRP 519	30.144	40.784	7.945	1.00	38.15	C2
ATOM 3573 CB TRP 519	31.124	40.083	8.780	1.00	38.52	C2
ATOM 3574 CG TRP 519	30.493	38.793	9.255	1.00	42.26	C2
ATOM 3575 CD2 TRP 519	29.880	38.578	10.473	1.00	41.70	C2
ATOM 3576 CE2 TRP 519	29.437	37.278	10.335	1.00	41.69	C2
ATOM 3577 CE3 TRP 519	29.648	39.282	11.629	1.00	42.26	C2
ATOM 3578 CD1 TRP 519	30.448	37.695	8.419	1.00	42.92	C2
ATOM 3579 NE1 TRP 519	29.788	36.793	9.115	1.00	44.19	C2
ATOM 3580 HE1 TRP 519	29.485	35.935	8.741	1.00	0.00	C2
ATOM 3581 C22 TRP 519	28.753	36.671	11.360	1.00	41.91	C2
ATOM 3582 C23 TRP 519	28.964	38.666	12.652	1.00	41.91	C2
ATOM 3583 CH2 TRP 519	28.522	37.375	12.515	1.00	41.05	C2
ATOM 3584 C TRP 519	29.027	41.368	8.815	1.00	39.33	C2
ATOM 3585 O TRP 519	27.888	40.919	8.726	1.00	38.28	C2
ATOM 3586 N GLN 520	29.264	42.375	9.650	1.00	41.86	C2
ATOM 3587 H GLN 520	30.180	42.717	9.700	1.00	0.00	C2
ATOM 3588 CA GLN 520	28.240	43.016	10.464	1.00	44.63	C2
ATOM 3589 CB GLN 520	28.691	44.198	11.239	1.00	47.03	C2
ATOM 3590 CG GLN 520	29.602	43.808	12.360	1.00	54.78	C2
ATOM 3591 CD GLN 520	29.910	45.009	13.243	1.00	60.14	C2
ATOM 3592 OE1 GLN 520	28.988	45.566	13.854	1.00	61.62	C2
ATOM 3593 NE1 GLN 520	31.172	45.456	13.371	1.00	60.46	C2
ATOM 3594 HE22 GLN 520	31.289	46.260	13.910	1.00	0.00	C2
ATOM 3595 IE22 GLN 520	31.895	44.966	12.932	1.00	0.00	C2
ATOM 3596 C GLN 520	27.141	43.577	9.621	1.00	46.28	C2
ATOM 3597 O GLN 520	26.001	43.474	10.059	1.00	48.62	C2
ATOM 3598 N GLN 521	27.362	44.145	8.442	1.00	46.99	C2
ATOM 3599 H GLN 521	28.272	44.257	8.092	1.00	0.00	C2
ATOM 3600 CA GLN 521	26.276	44.638	7.716	1.00	49.02	C2
ATOM 3601 CB GLN 521	26.632	45.533	6.566	1.00	50.06	C2
ATOM 3602 CG GLN 521	25.456	46.226	5.790	1.00	50.87	C2
ATOM 3603 CD GLN 521	24.616	47.278	6.534	1.00	51.82	C2
ATOM 3604 OE1 GLN 521	24.864	47.694	7.671	1.00	52.47	C2
ATOM 3605 NE2 GLN 521	23.577	47.776	5.888	1.00	50.36	C2
ATOM 3606 HE21 GLN 521	23.392	47.455	4.987	1.00	0.00	C2
ATOM 3607 HE22 GLN 521	23.044	48.424	6.390	1.00	0.00	C2
ATOM 3608 C GLN 521	25.454	43.446	7.155	1.00	50.15	C2
ATOM 3609 O GLN 521	24.214	43.514	7.177	1.00	51.82	C2
ATOM 3610 N MET 522	26.057	42.348	6.668	1.00	49.18	C2
ATOM 3611 H MET 522	27.038	42.291	6.688	1.00	0.00	C2
ATOM 3612 CA MET 522	25.280	41.227	6.171	1.00	48.22	C2
ATOM 3613 CB MET 522	26.185	40.167	5.607	1.00	46.36	C2
ATOM 3614 CG MET 522	26.942	40.661	4.412	1.00	44.32	C2
ATOM 3615 SD MET 522	27.855	39.435	3.426	1.00	48.35	C2
ATOM 3616 CE MET 522	28.795	38.447	4.565	1.00	47.80	C2
ATOM 3617 O MET 522	24.453	40.642	7.316	1.00	50.14	C2
ATOM 3618 C MET 522	23.380	40.124	7.038	1.00	50.60	C2
ATOM 3619 N GLU 523	24.848	40.772	8.596	1.00	51.91	C2
ATOM 3620 H GLU 523	25.766	41.031	8.769	1.00	0.00	C2
ATOM 3621 CA GLU 523	24.027	40.313	9.718	1.00	54.53	C2
ATOM 3622 CB GLU 523	24.654	40.486	11.081	1.00	54.40	C2
ATOM 3623 CG GLU 523	25.732	39.525	11.198	1.00	57.05	C2
ATOM 3624 CD GLU 523	25.386	38.150	10.888	1.00	61.72	C2
ATOM 3625 OE1 GLU 523	24.515	37.487	11.477	1.00	64.66	C2
ATOM 3626 OE2 GLU 523	25.979	37.773	9.872	1.00	63.19	C2
ATOM 3627 C GLU 523	22.773	41.116	9.536	1.00	58.29	C2
ATOM 3628 O GLU 523	21.688	40.538	9.850	1.00	59.32	C2
ATOM 3629 N ALA 524	22.920	42.432	9.992	1.00	61.41	C2
ATOM 3630 H ALA 524	23.814	42.798	10.024	1.00	0.00	C2
ATOM 3631 CA ALA 524	21.815	43.360	10.076	1.00	63.58	C2
ATOM 3632 CB ALA 524	22.382	44.768	9.992	1.00	64.11	C2
ATOM 3633 C ALA 524	20.818	43.109	8.946	1.00	64.79	C2
ATOM 3634 O ALA 524	19.655	42.824	9.206	1.00	65.69	C2
ATOM 3635 N ALA 525	21.251	43.083	7.693	1.00	66.44	C2
ATOM 3636 H ALA 525	22.196	43.283	7.516	1.00	68.58	C2
ATOM 3637 CA ALA 525	20.371	42.789	6.574	1.00	68.58	C2
ATOM 3638 CB ALA 525	21.117	43.044	5.288	1.00	67.42	C2
ATOM 3639 C ALA 525	19.841	41.356	6.558	1.00	71.11	C2
ATOM 3640 O ALA 525	19.116	40.946	5.651	1.00	71.65	C2
ATOM 3641 N GLY 526	20.257	40.510	7.498	1.00	74.20	C2
ATOM 3642 H GLY 526	21.019	40.780	8.043	1.00	0.00	C2
ATOM 3643 CA GLY 526	19.728	39.157	7.653	1.00	76.30	C2
ATOM 3644 C GLY 526	20.430	38.085	6.842	1.00	78.13	C2
ATOM 3645 O GLY 526	20.174	36.910	7.094	1.00	79.05	C2
ATOM 3646 N MET 527	21.388	38.433	5.970	1.00	80.23	C2
ATOM 3647 H MET 527	21.759	39.337	6.075	1.00	0.00	C2
ATOM 3648 CA MET 527	22.055	37.489	5.063	1.00	81.73	C2
ATOM 3649 CB MET 527	22.771	38.256	3.928	1.00	81.72	C2
ATOM 3650 CG MET 527	22.385	39.719	3.720	1.00	81.52	C2
ATOM 3651 SD MET 527	23.364	40.523	2.436	1.00	87.04	C2
ATOM 3652 CE MET 527	22.600	42.117	2.403	1.00	84.47	C2
ATOM 3653 C MET 527	73.078	36.584	5.780	1.00	82.04	C2
ATOM 3654 OT1 MET 527	22.974	35.357	5.624	1.00	83.38	C2
ATOM 3655 OT2 MET 527	23.949	37.104	6.500	1.00	82.99	C2
ATOM 3656 CB MET 538	47.224	28.531	2.401	1.00	77.43	C3
ATOM 3657 CG MET 538	47.397	30.041	2.427	1.00	77.15	C3
ATOM 3658 SD MET 538	46.205	30.708	3.604	1.00	79.03	C3
ATOM 3659 CE MET 538	48.850	31.067	2.515	1.00	77.20	C3
ATOM 3660 C MET 538	48.549	27.839	0.386	1.00	75.32	C3
ATOM 3661 O MET 538	49.130	26.745	0.405	1.00	77.11	C3
ATOM 3662 HT1 MET 538	47.563	26.068	1.449	1.00	0.00	C3
ATOM 3663 HT2 MET 538	46.638	26.204	0.075	1.00	0.00	C3
ATOM 3664 N MET 538	46.724	26.552	1.050	1.00	77.52	C3
ATOM 3665 HT3 MET 538	45.873	26.401	1.617	1.00	0.00	C3
ATOM 3666 CA MET 538	47.153	27.940	0.995	1.00	76.57	C3
ATOM 3667 H PRO 539	49.089	28.870	-0.224	1.00	77.65	C3
ATOM 3668 CD PRO 539	48.346	29.821	-1.046	1.00	72.26	C3
ATOM 3669 CA PRO 539	50.526	29.020	-0.549	1.00	72.34	C3
ATOM 3670 CB PRO 539	50.677	30.165	-1.106	1.00	71.49	C3

FIGURE 5

ATOM 3671	CG PRO	S39	-49.437	30.503	-1.837	1.00	71.52	C3	ATOM 3722	C	PIE	S45	61.543	34.900	-1.667	1.00	34.81	C3
ATOM 3672	C PRO	S39	51.250	28.931	0.991	1.00	67.83	C3	ATOM 3723	O	PIE	S45	60.901	35.660	-2.389	1.00	32.88	C3
ATOM 3673	O PRO	S39	50.666	29.294	2.091	1.00	68.05	C3	ATOM 3724	N	GLN	S46	60.912	33.847	-1.135	1.00	34.77	C3
ATOM 3674	N ALA	S40	52.484	28.417	0.961	1.00	64.48	C3	ATOM 3725	II	GLN	S46	61.396	33.223	-0.558	1.00	0.02	C3
ATOM 3675	II ALA	S40	52.858	28.098	0.111	1.00	0.00	C3	ATOM 3726	CA	GIN	S46	59.490	33.637	-1.433	1.00	33.72	C3
ATOM 3676	CA ALA	S40	53.389	28.498	2.112	1.00	61.83	C3	ATOM 3727	CG	GIN	S46	59.145	32.232	-1.140	1.00	34.85	C3
ATOM 3677	CB ALA	S40	54.004	27.200	2.619	1.00	63.57	C3	ATOM 3728	CG	GIN	S46	59.582	31.585	-2.444	1.00	42.45	C3
ATOM 3678	O ALA	S40	54.559	29.212	1.496	1.00	58.74	C3	ATOM 3729	CD	GIN	S46	59.374	30.085	-2.473	1.00	46.95	C3
ATOM 3679	O ALA	S40	54.835	29.036	0.301	1.00	58.30	C3	ATOM 3730	OII	GIN	S46	59.287	29.472	-1.399	1.00	48.90	C3
ATOM 3680	N PIE	S41	55.256	30.008	2.292	1.00	55.75	C3	ATOM 3731	NE2	GIN	S46	59.339	29.442	-3.644	1.00	47.20	C3
ATOM 3681	II PIE	S41	55.093	30.068	3.257	1.00	0.00	C3	ATOM 3732	II	GIN	S46	59.476	29.948	-4.472	1.00	0.00	C3
ATOM 3682	CA PIE	S41	56.299	30.814	1.702	1.00	51.38	C3	ATOM 3733	II	GIN	S46	59.154	28.481	-3.609	1.00	61.62	C3
ATOM 3683	CB PIE	S41	55.964	32.306	1.942	1.00	48.80	C3	ATOM 3734	C	GIN	S46	58.504	34.541	-0.729	1.00	61.62	C3
ATOM 3684	CG PIE	S41	54.789	32.703	1.058	1.00	45.77	C3	ATOM 3735	O	GIN	S46	57.429	34.850	-1.233	1.00	29.88	C3
ATOM 3685	CD1 PIE	S41	54.992	32.939	-0.279	1.00	44.20	C3	ATOM 3736	N	ARG	S47	58.907	34.929	0.465	1.00	31.09	C3
ATOM 3686	CD2 PIE	S41	53.507	32.747	1.582	1.00	44.76	C3	ATOM 3737	II	ARG	S47	59.750	34.566	0.811	1.00	0.90	C3
ATOM 3687	CE1 PIE	S41	53.901	32.207	-1.074	1.00	43.98	C3	ATOM 3738	CA	ARG	S47	58.160	35.830	1.282	1.00	31.43	C3
ATOM 3688	CE2 PIE	S41	52.428	33.018	0.769	1.00	42.86	C3	ATOM 3739	CB	ARG	S47	58.813	35.874	2.601	1.00	31.74	C3
ATOM 3689	CZ PIE	S41	52.625	33.247	-0.563	1.00	42.52	C3	ATOM 3740	CG	ARG	S47	57.906	35.224	3.623	1.00	37.02	C3
ATOM 3690	C PIE	S41	57.586	30.364	2.333	1.00	49.55	C3	ATOM 3741	CD	ARG	S47	58.344	33.858	4.016	1.00	40.56	C3
ATOM 3691	O PIE	S41	58.002	30.807	3.395	1.00	49.55	C3	ATOM 3742	NE	ARG	S47	59.743	34.058	4.345	1.00	47.90	C3
ATOM 3692	N ALA	S42	58.172	29.442	1.562	1.00	48.21	C3	ATOM 3743	IE	ARG	S47	60.389	33.924	3.620	1.00	0.00	C3
ATOM 3693	II ALA	S42	57.825	29.298	0.656	1.00	0.00	C3	ATOM 3744	CZ	ARG	S47	60.190	34.394	5.543	1.00	49.48	C3
ATOM 3694	CA ALA	S42	59.376	28.711	1.968	1.00	45.37	C3	ATOM 3745	NIH1	ARG	S47	59.361	34.522	6.593	1.00	51.97	C3
ATOM 3695	CB ALA	S42	59.700	27.749	0.898	1.00	45.21	C3	ATOM 3746	NIH11	ARG	S47	58.380	34.356	6.488	1.00	0.00	C3
ATOM 3696	C ALA	S42	60.510	29.567	2.766	1.00	44.87	C3	ATOM 3747	NIH12	ARG	S47	59.731	34.763	7.491	1.00	0.00	C3
ATOM 3697	O ALA	S42	61.001	29.504	3.374	1.00	46.49	C3	ATOM 3748	NIH2	ARG	S47	61.464	34.775	5.616	1.00	48.55	C3
ATOM 3698	N SER	S43	61.013	30.408	1.395	1.00	42.63	C3	ATOM 3749	NIH21	ARG	S47	62.025	34.803	4.788	1.00	0.00	C3
ATOM 3699	H SER	S43	60.477	30.685	0.630	1.00	0.00	C3	ATOM 3750	NIH22	ARG	S47	61.854	35.034	6.501	1.00	0.00	C3
ATOM 3700	CA SER	S43	62.253	31.108	1.708	1.00	40.31	C3	ATOM 3751	C	ARG	S47	58.167	37.181	0.590	1.00	32.76	C3
ATOM 3701	CB SER	S43	63.170	30.861	0.587	1.00	37.74	C3	ATOM 3752	O	ARG	S47	57.084	37.694	0.317	1.00	34.25	C3
ATOM 3702	CG SER	S43	62.391	31.181	-0.554	1.00	35.74	C3	ATOM 3753	N	ARG	S48	59.348	37.717	0.205	1.00	31.44	C3
ATOM 3703	HG SER	S43	61.824	30.423	-0.751	1.00	0.00	C3	ATOM 3754	CA	ARG	S48	60.148	37.203	0.444	1.00	0.00	C3
ATOM 3704	C SER	S43	62.087	32.613	1.896	1.00	40.88	C3	ATOM 3755	CA	ARG	S48	59.529	38.980	-0.555	1.00	30.01	C3
ATOM 3705	O SER	S43	61.016	33.115	1.536	1.00	42.63	C3	ATOM 3756	CB	ARG	S48	60.995	39.213	-0.949	1.00	25.42	C3
ATOM 3706	N ALA	S44	63.120	33.383	2.310	1.00	38.84	C3	ATOM 3757	CG	ARG	S48	61.820	39.361	0.494	1.00	26.11	C3
ATOM 3707	II ALA	S44	63.929	32.951	2.650	1.00	0.00	C3	ATOM 3758	CD	ARG	S48	63.280	39.158	-0.054	1.00	79.34	C3
ATOM 3708	CA ALA	S44	63.035	34.836	2.345	1.00	37.31	C3	ATOM 3759	NE	ARG	S48	64.044	39.162	1.189	1.00	32.36	C3
ATOM 3709	CB ALA	S44	64.340	35.450	2.808	1.00	35.74	C3	ATOM 3760	IE	ARG	S48	63.572	38.883	1.995	1.00	0.00	C3
ATOM 3710	C ALA	S44	62.723	35.372	0.947	1.00	37.06	C3	ATOM 3761	CZ	ARG	S48	65.344	39.518	1.325	1.00	32.66	C3
ATOM 3711	O ALA	S44	61.829	36.220	0.820	1.00	38.23	C3	ATOM 3762	NIH1	ARG	S48	66.159	39.923	0.335	1.00	34.98	C3
ATOM 3712	N PIE	S45	63.357	34.881	-0.130	1.00	35.72	C3	ATOM 3763	NIH11	ARG	S48	67.107	40.170	0.533	1.00	0.00	C3
ATOM 3713	II PIE	S45	64.131	34.298	-0.010	1.00	0.00	C3	ATOM 3764	NIH12	ARG	S48	65.812	39.981	-0.600	1.00	0.00	C3
ATOM 3714	CA PIE	S45	62.992	35.268	-1.484	1.00	33.66	C3	ATOM 3765	NIH2	ARG	S48	65.837	39.518	2.549	1.00	32.03	C3
ATOM 3715	CB PIE	S45	63.738	34.534	-2.593	1.00	29.71	C3	ATOM 3766	NIH21	ARG	S48	66.788	39.781	2.708	1.00	0.00	C3
ATOM 3716	CG PIE	S45	63.140	34.742	-3.990	1.00	29.62	C3	ATOM 3767	NIH22	ARG	S48	65.250	39.275	3.321	1.00	0.00	C3
ATOM 3717	CD1 PIE	S45	62.317	33.788	-4.557	1.00	29.80	C3	ATOM 3768	C	ARG	S48	58.713	38.997	-1.832	1.00	29.81	C3
ATOM 3718	CD2 PIE	S45	63.371	35.915	-4.689	1.00	31.90	C3	ATOM 3769	O	ARG	S48	57.778	39.790	-1.968	1.00	33.63	C3
ATOM 3719	CE1 PIE	S45	61.723	33.984	-5.795	1.00	28.61	C3	ATOM 3770	N	ALA	S49	58.979	38.102	-2.761	1.00	27.87	C3
ATOM 3720	CE2 PIE	S45	62.777	36.113	-5.928	1.00	31.00	C3	ATOM 3771	II	ALA	S49	59.684	37.436	-2.641	1.00	0.00	C3
ATOM 3721	CZ PIE	S45	61.955	35.150	-6.480	1.00	31.01	C3	ATOM 3772	CA	ALA	S49	58.227	38.045	-3.984	1.00	27.18	C3

FIGURE 5

ATOM 3773 CB ALA 549	58.797	36.934	-4.857	1.00	28.72	C3
ATOM 3774 C ALA 549	56.748	37.810	-3.770	1.00	25.91	C3
ATOM 3775 O ALA 549	55.896	38.337	-4.468	1.00	26.03	C3
ATOM 3776 N GLY 550	56.421	37.074	-2.748	1.00	26.53	C3
ATOM 3777 H GLY 550	57.103	36.657	-2.185	1.00	0.00	C3
ATOM 3778 CA GLY 550	55.055	36.805	-2.457	1.00	26.08	C3
ATOM 3779 C GLY 550	54.410	38.098	-2.075	1.00	26.94	C3
ATOM 3780 O GLY 550	53.339	38.380	-2.608	1.00	26.59	C3
ATOM 3781 N GLY 551	55.073	38.917	-1.234	1.00	27.78	C3
ATOM 3782 H GLY 551	55.958	38.642	-0.925	1.00	0.00	C3
ATOM 3783 CA GLY 551	54.540	40.212	-0.779	1.00	26.51	C3
ATOM 3784 C GLY 551	54.302	41.113	-1.994	1.00	26.82	C3
ATOM 3785 O GLY 551	53.313	41.852	-2.065	1.00	27.82	C3
ATOM 3786 N VAL 552	55.154	41.013	-3.012	1.00	25.81	C3
ATOM 3787 H VAL 552	55.916	40.396	-2.954	1.00	0.00	C3
ATOM 3788 CA VAL 552	54.952	41.843	-4.176	1.00	28.39	C3
ATOM 3789 C VAL 552	56.178	41.743	-5.190	1.00	26.20	C3
ATOM 3790 CG1 VAL 552	55.917	42.391	-6.541	1.00	26.53	C3
ATOM 3791 CG2 VAL 552	57.327	42.546	-4.594	1.00	26.44	C3
ATOM 3792 C VAL 552	53.650	41.406	-4.820	1.00	29.05	C3
ATOM 3793 O VAL 552	52.744	42.251	-4.888	1.00	31.68	C3
ATOM 3794 N LEU 553	53.455	40.120	-5.176	1.00	27.20	C3
ATOM 3795 H LEU 553	54.122	39.447	-4.908	1.00	0.00	C3
ATOM 3796 CA LEU 553	52.266	39.705	-5.915	1.00	23.80	C3
ATOM 3797 C LEU 553	52.357	38.262	-6.363	1.00	24.86	C3
ATOM 3798 CG LEU 553	53.431	37.955	-7.357	1.00	23.06	C3
ATOM 3799 CD1 LEU 553	54.073	36.623	-7.092	1.00	24.31	C3
ATOM 3800 CD2 LEU 553	52.794	38.061	-8.703	1.00	21.87	C3
ATOM 3801 C LEU 553	51.012	39.825	-5.114	1.00	23.72	C3
ATOM 3802 O LEU 553	49.982	40.138	-5.712	1.00	24.63	C3
ATOM 3803 N VAL 554	50.962	39.580	-3.803	1.00	24.37	C3
ATOM 3804 H VAL 554	51.774	39.350	-3.295	1.00	0.00	C3
ATOM 3805 CA VAL 554	49.660	39.691	-3.180	1.00	26.36	C3
ATOM 3806 C VAL 554	49.472	38.751	-1.802	1.00	26.55	C3
ATOM 3807 CG1 VAL 554	50.696	37.933	-1.418	1.00	23.95	C3
ATOM 3808 CG2 VAL 554	48.953	39.614	-0.682	1.00	25.58	C3
ATOM 3809 C VAL 554	49.322	41.175	-2.960	1.00	27.53	C3
ATOM 3810 O VAL 554	48.142	41.502	-3.192	1.00	27.44	C3
ATOM 3811 N ALA 555	50.277	42.106	-2.716	1.00	28.04	C3
ATOM 3812 H ALA 555	51.221	41.831	-2.658	1.00	0.00	C3
ATOM 3813 CA ALA 555	49.956	43.539	-2.509	1.00	28.57	C3
ATOM 3814 CB ALA 555	51.161	44.427	-2.217	1.00	28.07	C3
ATOM 3815 C ALA 555	49.402	44.055	-3.803	1.00	28.12	C3
ATOM 3816 O ALA 555	48.425	44.803	-3.847	1.00	30.12	C3
ATOM 3817 N SER 556	49.985	43.521	-4.839	1.00	26.44	C3
ATOM 3818 H SER 556	50.781	42.956	-4.710	1.00	0.00	C3
ATOM 3819 CA SER 556	49.548	43.810	-6.152	1.00	30.09	C3
ATOM 3820 CB SER 556	50.684	43.277	-6.965	1.00	31.42	C3
ATOM 3821 OG SER 556	50.442	43.338	-8.344	1.00	37.88	C3
ATOM 3822 HG SER 556	49.966	44.144	-8.576	1.00	0.00	C3
ATOM 3823 C SER 556	48.143	43.243	-6.454	1.00	32.78	C3
ATOM 3824 O SER 556	47.287	43.961	-7.003	1.00	34.56	C3
ATOM 3825 N HIS 557	47.750	42.019	-6.088	1.00	32.78	C3
ATOM 3826 H HIS 557	48.350	41.453	-5.560	1.00	0.00	C3
ATOM 3827 CA HIS 557	46.396	41.605	-6.401	1.00	33.64	C3
ATOM 3828 CB HIS 557	46.203	40.142	-6.242	1.00	37.88	C3
ATOM 3829 CG HIS 557	46.986	39.518	-7.348	1.00	42.44	C3
ATOM 3830 CD2 HIS 557	46.694	39.665	-8.675	1.00	43.63	C3
ATOM 3831 ND1 HIS 557	48.108	38.837	-7.209	1.00	45.23	C3
ATOM 3832 HD1 HIS 557	48.641	38.764	-6.385	1.00	0.00	C3
ATOM 3833 CE1 HIS 557	48.524	38.569	-8.414	1.00	46.56	C3
ATOM 3834 NE2 HIS 557	47.676	39.066	-9.283	1.00	45.62	C3
ATOM 3835 HE2 HIS 557	47.793	39.018	-10.257	1.00	0.00	C3
ATOM 3836 C HIS 557	45.383	42.249	-5.520	1.00	32.94	C3
ATOM 3837 O HIS 557	44.256	42.444	-5.934	1.00	33.08	C3
ATOM 3838 N LEU 558	45.744	42.534	-4.280	1.00	33.05	C3
ATOM 3839 H LEU 558	46.657	42.356	-3.986	1.00	0.00	C3
ATOM 3840 CA LEU 558	44.817	43.125	-3.348	1.00	31.91	C3
ATOM 3841 CB LEU 558	45.420	43.107	-1.965	1.00	29.25	C3
ATOM 3842 CG LEU 558	44.605	43.615	-0.818	1.00	26.02	C3
ATOM 3843 CD1 LEU 558	43.279	42.883	-0.742	1.00	25.00	C3
ATOM 3844 CD2 LEU 558	45.496	43.571	0.408	1.00	22.86	C3
ATOM 3845 C LEU 558	44.527	44.521	-3.783	1.00	32.47	C3
ATOM 3846 O LEU 558	43.402	44.944	-3.596	1.00	33.97	C3
ATOM 3847 N GLN 559	45.482	45.231	-4.370	1.00	34.36	C3
ATOM 3848 H GLN 559	46.386	44.855	-4.406	1.00	0.00	C3
ATOM 3849 CA GLN 559	45.255	46.569	-4.912	1.00	36.75	C3
ATOM 3850 CB GLN 559	46.598	47.067	-5.470	1.00	39.63	C3
ATOM 3851 CG GLN 559	46.707	48.543	-5.875	1.00	42.07	C3
ATOM 3852 CD GLN 559	46.530	49.618	-4.793	1.00	41.53	C3
ATOM 3853 OE1 GLN 559	45.961	50.687	-5.057	1.00	41.25	C3
ATOM 3854 NE2 GLN 559	46.951	49.416	-3.561	1.00	37.96	C3
ATOM 3855 HE2 GLN 559	47.271	48.528	-3.311	1.00	0.00	C3
ATOM 3856 HE22 GLN 559	47.001	50.190	-2.967	1.00	0.00	C3
ATOM 3857 C GLN 559	44.142	46.635	-5.976	1.00	35.22	C3
ATOM 3858 O GLN 559	43.165	47.404	-5.839	1.00	34.99	C3
ATOM 3859 N SER 560	44.260	45.817	-7.025	1.00	33.46	C3
ATOM 3860 H SER 560	45.083	45.292	-7.154	1.00	0.00	C3
ATOM 3861 CA SER 560	43.222	45.683	-8.049	1.00	32.88	C3
ATOM 3862 CB SER 560	43.693	44.776	-9.088	1.00	34.05	C3
ATOM 3863 OG SER 560	45.021	45.174	-9.281	1.00	42.40	C3
ATOM 3864 HG SER 560	45.042	45.996	-9.783	1.00	0.00	C3
ATOM 3865 C SER 560	41.885	45.133	-7.559	1.00	32.01	C3
ATOM 3866 O SER 560	40.791	45.582	-7.920	1.00	32.23	C3
ATOM 3867 H PHE 561	41.969	44.123	-6.710	1.00	29.50	C3
ATOM 3868 H PHE 561	42.850	43.767	-6.464	1.00	0.00	C3
ATOM 3869 CA PHE 561	40.803	43.529	-6.118	1.00	28.37	C3
ATOM 3870 CB PHE 561	41.237	42.541	-5.040	1.00	26.27	C3
ATOM 3871 CG PHE 561	40.069	41.966	-4.268	1.00	25.68	C3
ATOM 3872 CD1 PHE 561	39.282	40.999	-4.846	1.00	25.44	C3
ATOM 3873 CD2 PHE 561	39.761	42.482	-3.051	1.00	25.45	C3
ATOM 3874 CE1 PHE 561	38.166	40.551	-4.215	1.00	21.49	C3

FIGURES

ATOM 3875	CE2 PHE	S61	38.635	42.027	-2.421	1.00	26.89	C3	ATOM 3926	C TYR	S66	34.534	50.696	-6.217	1.00	48.48	C3
ATOM 3876	C2 PHE	S61	37.853	41.074	-3.008	1.00	24.29	C3	ATOM 3927	O TYR	S66	33.545	51.376	-5.950	1.00	46.35	C3
ATOM 3877	C PHE	S61	39.987	44.645	-5.505	1.00	28.81	C3	ATOM 3928	N ALA	S67	34.679	50.115	-7.417	1.00	49.14	C3
ATOM 3878	O PHE	S61	38.789	44.697	-5.731	1.00	29.31	C3	ATOM 3929	H ALA	S67	35.512	49.625	-7.572	1.00	0.00	C3
ATOM 3879	N LEU	S62	40.672	45.565	-4.797	1.00	28.39	C3	ATOM 3930	CA ALA	S67	33.670	50.165	-8.490	1.00	52.09	C3
ATOM 3880	H LEU	S62	41.643	45.462	-4.707	1.00	0.00	C3	ATOM 3931	CB ALA	S67	34.210	49.574	-9.788	1.00	48.37	C3
ATOM 3881	CA LEU	S62	40.033	46.617	-4.057	1.00	26.51	C3	ATOM 3932	C ALA	S67	32.315	49.449	-8.248	1.00	55.31	C3
ATOM 3882	CB LEU	S62	40.964	47.203	-3.074	1.00	23.80	C3	ATOM 3933	O ALA	S67	31.276	50.008	-8.501	1.00	56.87	C3
ATOM 3883	CG LEU	S62	41.047	46.411	-1.816	1.00	24.85	C3	ATOM 3934	N VAL	S68	32.247	48.211	-7.736	1.00	57.66	C3
ATOM 3884	CD1 LEU	S62	42.207	46.868	-1.049	1.00	24.07	C3	ATOM 3935	H VAL	S68	33.083	47.729	-7.564	1.00	0.00	C3
ATOM 3885	CD2 LEU	S62	39.794	46.551	-1.008	1.00	25.13	C3	ATOM 3936	CA VAL	S68	31.119	46.031	-7.339	1.00	58.96	C3
ATOM 3886	C LEU	S62	39.586	47.669	-4.988	1.00	29.27	C3	ATOM 3937	CB VAL	S68	31.239	45.508	-5.911	1.00	60.27	C3
ATOM 3887	O LEU	S62	38.580	48.304	-4.681	1.00	29.83	C3	ATOM 3938	CG1 VAL	S68	29.851	45.471	-7.922	1.00	60.44	C3
ATOM 3888	N GLU	S63	40.239	47.871	-6.115	1.00	30.88	C3	ATOM 3939	CG2 VAL	S68	30.393	48.177	-6.245	1.00	62.66	C3
ATOM 3889	H GLU	S63	41.052	47.368	-6.325	1.00	0.00	C3	ATOM 3940	C VAL	S68	31.075	48.154	-6.180	1.00	64.78	C3
ATOM 3890	CA GLU	S63	39.738	48.908	-6.966	1.00	36.88	C3	ATOM 3941	O VAL	S68	32.847	49.687	-1.140	1.00	70.17	C3
ATOM 3891	CB GLU	S63	40.660	49.142	-8.137	1.00	40.80	C3	ATOM 3942	N LEU	S69	31.039	48.054	-1.286	1.00	70.56	C3
ATOM 3892	CG GLU	S63	41.999	49.628	-7.682	1.00	48.55	C3	ATOM 3943	H LEU	S69	29.567	50.509	-4.667	1.00	72.69	C3
ATOM 3893	CD GLU	S63	43.148	49.277	-8.619	1.00	55.42	C3	ATOM 3944	CA LEU	S69	28.365	50.553	-4.425	1.00	73.80	C3
ATOM 3894	OE1 GLU	S63	44.301	49.283	-8.135	1.00	57.39	C3	ATOM 3945	CB LEU	S69	30.180	51.391	-5.479	1.00	75.95	C3
ATOM 3895	OE2 GLU	S63	42.886	48.986	-9.808	1.00	56.44	C3	ATOM 3946	CG LEU	S69	31.553	51.299	-5.580	1.00	0.00	C3
ATOM 3896	C GLU	S63	38.375	48.469	-7.466	1.00	39.02	C3	ATOM 3947	CD1 LEU	S69	29.510	52.498	-6.173	1.00	78.78	C3
ATOM 3897	O GLU	S63	37.388	49.170	-7.270	1.00	39.09	C3	ATOM 3948	CD2 LEU	S69	30.399	53.068	-7.308	1.00	80.07	C3
ATOM 3898	N VAL	S64	38.289	47.255	-8.030	1.00	42.30	C3	ATOM 3949	C LEU	S69	29.658	54.222	-7.997	1.00	84.16	C3
ATOM 3899	H VAL	S64	39.107	46.714	-8.074	1.00	0.00	C3	ATOM 3950	O LEU	S69	29.976	54.744	-9.417	1.00	85.66	C3
ATOM 3900	CA VAL	S64	37.052	46.683	-8.558	1.00	41.84	C3	ATOM 3951	N ARG	S70	28.892	55.690	-9.737	1.00	85.67	C3
ATOM 3901	CB VAL	S64	37.333	45.255	-9.041	1.00	42.27	C3	ATOM 3952	H ARG	S70	27.971	55.354	-9.727	1.00	0.00	C3
ATOM 3902	CG1 VAL	S64	36.055	44.538	-9.435	1.00	41.17	C3	ATOM 3953	CA ARG	S70	29.051	56.991	-10.026	1.00	85.06	C3
ATOM 3903	CG2 VAL	S64	38.283	45.348	-10.241	1.00	42.11	C3	ATOM 3954	CB ARG	S70	30.240	57.590	-10.082	1.00	84.43	C3
ATOM 3904	C VAL	S64	36.030	46.709	-7.442	1.00	41.68	C3	ATOM 3955	CG ARG	S70	31.069	57.036	-9.908	1.00	0.00	C3
ATOM 3905	O VAL	S64	34.892	47.015	-7.697	1.00	42.34	C3	ATOM 3956	CD ARG	S70	30.295	58.561	-10.314	1.00	0.00	C3
ATOM 3906	N SER	S65	36.419	46.501	-6.706	1.00	42.75	C3	ATOM 3957	NE ARG	S70	27.958	57.736	-10.154	1.00	84.57	C3
ATOM 3907	H SER	S65	37.333	46.173	-6.063	1.00	0.00	C3	ATOM 3958	HE ARG	S70	27.059	57.316	-10.030	1.00	0.00	C3
ATOM 3908	CA SER	S65	35.562	46.602	-5.064	1.00	44.85	C3	ATOM 3959	C2 ARG	S70	28.042	58.708	-10.375	1.00	0.00	C3
ATOM 3909	CB SER	S65	36.344	46.013	-3.894	1.00	46.54	C3	ATOM 3960	NH1 ARG	S70	28.201	52.009	-6.812	1.00	79.92	C3
ATOM 3910	OG SER	S65	35.590	45.714	-2.731	1.00	51.75	C3	ATOM 3961	HIH1 ARG	S70	27.107	52.565	-6.709	1.00	79.61	C3
ATOM 3911	HG SER	S65	35.060	46.481	-2.491	1.00	0.00	C3	ATOM 3962	HIH2 ARG	S70	28.362	50.900	-7.511	1.00	81.35	C3
ATOM 3912	C SER	S65	35.167	48.063	-4.871	1.00	45.70	C3	ATOM 3963	HIH2 ARG	S70	29.214	50.417	-7.440	1.00	0.00	C3
ATOM 3913	O SER	S65	34.038	48.287	-4.446	1.00	46.87	C3	ATOM 3964	HIH2 ARG	S70	27.247	50.106	-8.197	1.00	82.75	C3
ATOM 3914	N TYR	S66	35.965	49.093	-5.146	1.00	47.59	C3	ATOM 3965	HIH2 ARG	S70	27.882	49.274	-9.167	1.00	83.42	C3
ATOM 3915	H TYR	S66	36.893	48.908	-5.386	1.00	0.00	C3	ATOM 3966	C ARG	S70	28.633	50.079	-10.280	1.00	85.08	C3
ATOM 3916	CA TYR	S66	35.518	50.474	-5.086	1.00	49.68	C3	ATOM 3967	O ARG	S70	28.971	49.529	-11.532	1.00	85.81	C3
ATOM 3917	CB TYR	S66	36.765	51.362	-5.164	1.00	56.17	C3	ATOM 3968	N IIS	S71	29.074	51.303	-10.268	1.00	86.25	C3
ATOM 3918	CG TYR	S66	36.715	52.622	-6.007	1.00	64.64	C3	ATOM 3969	H IIS	S71	29.080	51.900	-9.489	1.00	0.00	C3
ATOM 3919	CD1 TYR	S66	37.264	52.538	-7.278	1.00	69.53	C3	ATOM 3970	CA IIS	S71	29.595	51.595	-11.439	1.00	86.31	C3
ATOM 3920	CE1 TYR	S66	37.212	53.613	-8.151	1.00	73.71	C3	ATOM 3971	CB IIS	S71						
ATOM 3921	CD2 TYR	S66	36.109	53.797	-5.569	1.00	67.98	C3	ATOM 3972	CG IIS	S71						
ATOM 3922	CE2 TYR	S66	36.048	54.888	-6.441	1.00	72.92	C3	ATOM 3973	CD2 IIS	S71						
ATOM 3923	C2 TYR	S66	36.599	54.787	-7.735	1.00	75.29	C3	ATOM 3974	NH1 IIS	S71						
ATOM 3924	OH TYR	S66	36.538	55.838	-8.652	1.00	77.42	C3	ATOM 3975	HIH1 IIS	S71						
ATOM 3925	IH1 TYR	S66	36.905	55.565	-9.494	1.00	0.00	C3	ATOM 3976	CE1 IIS	S71						

FIGURE 5

ATOM 3977 NE2 H1S 571	29.494	50.518	-12.187	1.00	86.28	C3	ATOM 4028 H1 H2O 622	24.393	32.417	14.215	1.00	0.00	W
ATOM 3978 HE2 H1S 571	29.801	50.468	-13.119	1.00	0.00	C3	ATOM 4029 H2 H2O 622	24.469	31.478	13.112	1.00	0.00	W
ATOM 3979 C H1S 571	26.225	49.759	-7.195	1.00	83.31	C3	ATOM 4030 OH2 H2O 623	20.791	28.583	14.218	1.00	50.17	W
ATOM 3980 O H1S 571	25.075	50.194	-7.301	1.00	84.06	C3	ATOM 4031 H1 H2O 623	20.499	28.803	13.375	1.00	0.00	W
ATOM 3981 N LEU 572	26.540	48.963	-6.158	1.00	83.11	C3	ATOM 4032 H2 H2O 623	19.939	28.549	14.688	1.00	0.00	W
ATOM 3982 H LEU 572	27.474	48.824	-5.915	1.00	83.71	C3	ATOM 4033 OH2 H2O 625	22.680	78.881	2.761	1.00	40.48	W
ATOM 3983 CA LEU 572	25.527	48.457	-5.241	1.00	83.71	C3	ATOM 4034 H1 H2O 625	21.938	78.856	3.375	1.00	0.00	W
ATOM 3984 CB LEU 572	26.085	47.267	-4.454	1.00	83.57	C3	ATOM 4035 H2 H2O 625	22.766	79.246	1.970	1.00	0.00	W
ATOM 3985 CG LEU 572	25.439	45.884	-4.721	1.00	83.79	C3	ATOM 4036 OH2 H2O 626	39.689	36.486	9.730	1.00	23.46	W
ATOM 3986 CD1 LEU 572	25.783	45.386	-6.127	1.00	84.16	C3	ATOM 4037 H1 H2O 626	39.090	35.724	9.672	1.00	0.00	W
ATOM 3987 CD2 LEU 572	25.958	44.866	-3.714	1.00	84.08	C3	ATOM 4038 H2 H2O 626	39.627	36.872	8.853	1.00	0.00	W
ATOM 3988 C LEU 572	24.997	49.511	-4.261	1.00	84.78	C3	ATOM 4039 OH2 H2O 627	42.035	78.320	5.697	1.00	46.19	W
ATOM 3989 O LEU 572	24.265	49.192	-3.295	1.00	84.85	C3	ATOM 4040 H1 H2O 627	42.416	77.450	5.832	1.00	0.00	W
ATOM 3990 N ALA 573	25.349	50.796	-4.483	1.00	85.56	C3	ATOM 4041 H2 H2O 627	41.243	78.146	5.181	1.00	0.00	W
ATOM 3991 H ALA 573	26.020	50.980	-5.174	1.00	0.00	C3	ATOM 4042 OH2 H2O 631	47.222	31.440	6.299	1.00	34.17	W
ATOM 3992 CA ALA 573	24.822	51.925	-3.721	1.00	85.90	C3	ATOM 4043 H1 H2O 631	47.533	32.209	5.809	1.00	0.00	W
ATOM 3993 CB ALA 573	25.600	53.207	-3.970	1.00	85.79	C3	ATOM 4044 H2 H2O 631	47.442	30.713	5.714	1.00	0.00	W
ATOM 3994 C ALA 573	23.373	52.245	-4.057	1.00	87.21	C3	ATOM 4045 OH2 H2O 636	24.043	65.423	-0.336	1.00	73.38	W
ATOM 3995 OT1 ALA 573	22.610	52.413	-3.099	1.00	88.33	C3	ATOM 4046 H1 H2O 636	24.179	65.781	-1.228	1.00	0.00	W
ATOM 3996 OT2 ALA 573	23.022	52.309	-5.248	1.00	88.34	C3	ATOM 4047 H2 H2O 636	23.469	66.096	0.054	1.00	0.00	W
ATOM 3997 OH2 H2O 603	26.735	24.280	5.161	1.00	27.42	W	ATOM 4048 OH2 H2O 638	38.984	67.955	-11.226	1.00	29.97	W
ATOM 3998 H1 H2O 603	27.332	24.335	4.407	1.00	0.00	W	ATOM 4049 H1 H2O 638	38.283	67.402	-11.580	1.00	0.00	W
ATOM 3999 H2 H2O 603	26.288	23.435	4.992	1.00	0.00	W	ATOM 4050 H2 H2O 638	39.568	68.046	-11.998	1.00	0.00	W
ATOM 4000 OH2 H2O 605	47.880	37.960	12.023	1.00	56.30	W	ATOM 4051 OH2 H2O 639	27.930	66.675	-7.733	1.00	43.40	W
ATOM 4001 H1 H2O 605	47.789	37.874	13.031	1.00	0.00	W	ATOM 4052 H1 H2O 639	28.192	67.028	-6.876	1.00	0.00	W
ATOM 4002 H2 H2O 605	46.980	37.858	11.753	1.00	0.00	W	ATOM 4053 H2 H2O 639	26.975	66.791	-7.705	1.00	0.00	W
ATOM 4003 OH2 H2O 607	40.001	49.224	7.214	1.00	40.04	W	ATOM 4054 OH2 H2O 643	50.619	62.802	0.813	1.00	36.55	W
ATOM 4004 H1 H2O 607	40.471	48.761	7.909	1.00	0.00	W	ATOM 4055 H1 H2O 643	51.575	62.904	0.824	1.00	0.00	W
ATOM 4005 H2 H2O 607	40.123	48.642	6.457	1.00	0.00	W	ATOM 4056 H2 H2O 643	50.301	63.665	0.525	1.00	0.00	W
ATOM 4006 OH2 H2O 610	59.883	42.530	-9.698	1.00	38.90	W	ATOM 4057 OH2 H2O 646	62.897	38.367	3.759	1.00	73.55	W
ATOM 4007 H1 H2O 610	60.512	41.833	-9.477	1.00	0.00	W	ATOM 4058 H1 H2O 646	62.414	38.098	2.978	1.00	0.00	W
ATOM 4008 H2 H2O 610	59.189	42.046	-10.160	1.00	0.00	W	ATOM 4059 H2 H2O 646	62.244	38.247	4.461	1.00	0.00	W
ATOM 4009 OH2 H2O 611	57.178	35.940	-14.220	1.00	34.63	W	ATOM 4060 OH2 H2O 650	29.587	68.480	-9.555	1.00	65.67	W
ATOM 4010 H1 H2O 611	57.174	36.545	-14.974	1.00	0.00	W	ATOM 4061 H1 H2O 650	28.846	68.630	-10.148	1.00	0.00	W
ATOM 4011 H2 H2O 611	57.989	36.211	-13.757	1.00	0.00	W	ATOM 4062 H2 H2O 650	29.180	67.844	-8.936	1.00	0.00	W
ATOM 4012 OH2 H2O 612	25.793	27.337	19.130	1.00	29.21	W	ATOM 4063 OH2 H2O 652	51.408	56.331	-4.056	1.00	62.90	W
ATOM 4013 H1 H2O 612	26.709	27.661	19.145	1.00	0.00	W	ATOM 4064 H1 H2O 652	50.718	56.353	3.365	1.00	0.00	W
ATOM 4014 H2 H2O 612	25.762	26.792	19.929	1.00	0.00	W	ATOM 4065 H2 H2O 652	51.052	55.671	4.648	1.00	0.00	W
ATOM 4015 OH2 H2O 615	29.766	34.284	9.444	1.00	45.03	W	ATOM 4066 OH2 H2O 653	49.404	56.072	2.161	1.00	51.28	W
ATOM 4016 H1 H2O 615	30.017	34.618	10.308	1.00	0.00	W	ATOM 4067 H1 H2O 653	49.442	55.351	1.474	1.00	0.00	W
ATOM 4017 H2 H2O 615	29.113	33.592	9.660	1.00	0.00	W	ATOM 4068 H2 H2O 653	49.323	56.829	1.630	1.00	0.00	W
ATOM 4018 OH2 H2O 617	37.316	40.012	10.872	1.00	35.21	W	ATOM 4069 OH2 H2O 654	68.215	42.294	-2.563	1.00	40.77	W
ATOM 4019 H1 H2O 617	36.600	40.017	11.519	1.00	0.00	W	ATOM 4070 H1 H2O 654	68.347	41.745	-1.777	1.00	0.00	W
ATOM 4020 H2 H2O 617	37.944	39.376	11.259	1.00	0.00	W	ATOM 4071 H2 H2O 654	68.189	43.181	-2.190	1.00	0.00	W
ATOM 4021 OH2 H2O 619	40.370	52.041	-7.387	1.00	29.62	W	ATOM 4072 OH2 H2O 655	66.374	40.425	-2.489	1.00	47.31	W
ATOM 4022 H1 H2O 619	40.672	52.724	-6.779	1.00	0.00	W	ATOM 4073 H1 H2O 655	66.936	41.162	-2.766	1.00	0.00	W
ATOM 4023 H2 H2O 619	39.505	51.810	-7.052	1.00	0.00	W	ATOM 4074 H2 H2O 655	66.452	39.841	-3.252	1.00	0.00	W
ATOM 4024 OH2 H2O 621	27.903	32.440	10.664	1.00	39.99	W	ATOM 4075 OH2 H2O 656	66.927	41.478	-5.011	1.00	44.08	W
ATOM 4025 H1 H2O 621	27.553	33.207	11.141	1.00	0.00	W	ATOM 4076 H1 H2O 656	66.207	42.071	-4.989	1.00	0.00	W
ATOM 4026 H2 H2O 621	27.929	31.808	11.398	1.00	0.00	W	ATOM 4077 H2 H2O 656	67.542	41.824	-4.374	1.00	0.00	W
ATOM 4027 OH2 H2O 622	25.057	31.972	13.675	1.00	32.70	W	ATOM 4078 OH2 H2O 657	40.371	57.111	5.730	1.00	66.55	W

FIGURE 5

ATOM	4079	H1	H2O	657	39.958	56.259	5.613	1.00	0.00	W
ATOM	4080	H2	H2O	657	40.021	57.651	5.014	1.00	0.00	W
ATOM	4081	OH2	H2O	658	48.780	47.580	-3.122	1.00	52.09	W
ATOM	4082	H1	H2O	658	48.811	46.671	-3.438	1.00	0.00	W
ATOM	4083	H2	H2O	658	49.568	47.955	-3.542	1.00	0.00	W
ATOM	4084	OH2	H2O	663	29.095	62.889	1.825	1.00	39.23	W
ATOM	4085	H1	H2O	663	29.380	62.827	2.739	1.00	0.00	W
ATOM	4086	H2	H2O	663	28.377	63.526	1.887	1.00	0.00	W
ATOM	4087	OH2	H2O	664	27.132	75.640	7.430	1.00	50.65	W
ATOM	4088	H1	H2O	664	26.870	24.838	7.876	1.00	0.00	W
ATOM	4089	H2	H2O	664	27.001	25.362	6.496	1.00	0.00	W
ATOM	4090	OH2	H2O	665	23.367	30.554	12.167	1.00	49.69	W
ATOM	4091	H1	H2O	665	24.026	30.006	11.707	1.00	0.00	W
ATOM	4092	H2	H2O	665	22.941	31.016	11.438	1.00	0.00	W
ATOM	4093	OH2	H2O	666	46.015	32.192	10.179	1.00	66.86	W
ATOM	4094	H1	H2O	666	46.060	31.519	9.497	1.00	0.00	W
ATOM	4095	H2	H2O	666	45.411	31.827	10.833	1.00	0.00	W
ATOM	4096	OH2	H2O	667	38.943	37.883	11.978	1.00	47.87	W
ATOM	4097	H1	H2O	667	39.367	37.487	11.188	1.00	0.00	W
ATOM	4098	H2	H2O	667	38.521	37.114	12.362	1.00	0.00	W
ATOM	4099	OH2	H2O	671	33.437	58.101	2.269	1.00	46.65	W
ATOM	4100	H1	H2O	671	33.555	57.162	2.433	1.00	0.00	W
ATOM	4101	H2	H2O	671	33.962	58.514	2.961	1.00	0.00	W
ATOM	4102	OH2	H2O	672	27.551	31.314	20.022	1.00	30.15	W
ATOM	4103	H1	H2O	672	27.929	32.042	20.533	1.00	0.00	W
ATOM	4104	H2	H2O	672	26.845	31.764	19.552	1.00	0.00	W
ATOM	4105	OH2	H2O	673	25.714	36.908	21.385	1.00	36.95	W
ATOM	4106	H1	H2O	673	24.806	37.123	21.637	1.00	0.00	W
ATOM	4107	H2	H2O	673	25.599	36.284	20.654	1.00	0.00	W
ATOM	4108	OH2	H2O	674	38.244	66.897	12.076	1.00	57.36	W
ATOM	4109	H1	H2O	674	37.773	67.536	12.626	1.00	0.00	W
ATOM	4110	H2	H2O	674	38.153	66.104	12.618	1.00	0.00	W
ATOM	4111	OH2	H2O	675	35.762	36.553	-3.986	1.00	58.40	W
ATOM	4112	H1	H2O	675	35.600	37.449	-3.677	1.00	0.00	W
ATOM	4113	H2	H2O	675	35.549	36.642	-4.923	1.00	0.00	W
ATOM	4114	OH2	H2O	676	30.689	32.814	25.675	1.00	59.30	W
ATOM	4115	H1	H2O	676	30.093	33.571	25.680	1.00	0.00	W
ATOM	4116	H2	H2O	676	31.550	33.214	25.540	1.00	0.00	W
END										

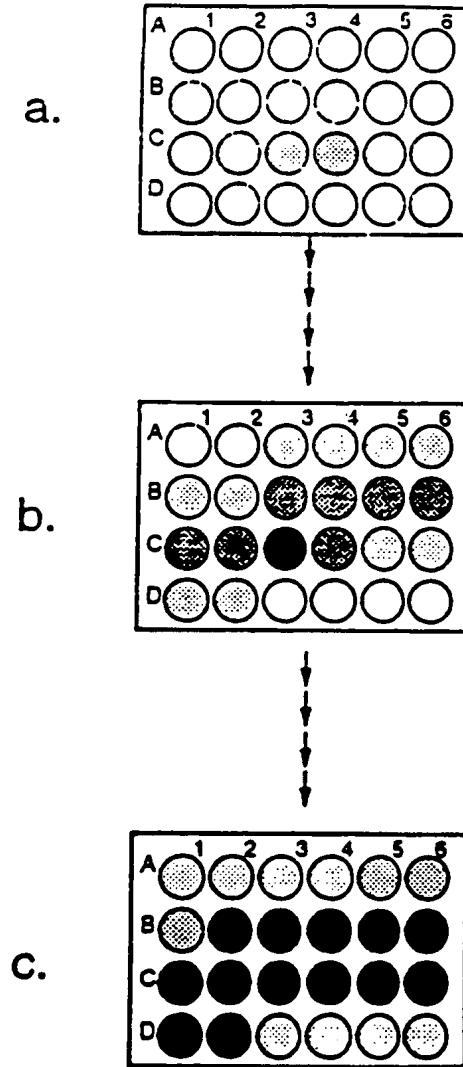


FIGURE 6



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Application Number
EP 94 10 1207

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The present search report has been drawn up for all claims			
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CATEGORY OF CITED DOCUMENTS X : particularly relevant if taken alone Y : particularly relevant if combined with another document of the same category A : technological background O : non-written disclosure P : intermediate document T : theory or principle underlying the invention E : earlier patent document, but published on, or after the filing date D : document cited in the application L : document cited for other reasons @ : member of the same patent family, corresponding document			



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The present search report has been drawn up for all claims			
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